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U.S. Army Center for Health Promotion
and Preventive Medicine

**TRAINING MUNITIONS HEALTH RISK
ASSESSMENT
NO. 39-EJ-1485-00
RESIDENTIAL EXPOSURE FROM INHALATION OF
AIR EMISSIONS FROM THE
LONG RIFLE .22 CALIBER BALL CARTRIDGE
DEPARTMENT OF DEFENSE IDENTIFICATION CODE: A106**

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Readiness Thru Health

U.S. Army Center for Health Promotion and Preventive Medicine

The lineage of the U.S. Army Center for Health Promotion and Preventive Medicine (USACHPPM) can be traced back over 50 years. This organization began as the U.S. Army Industrial Hygiene Laboratory, established during the industrial buildup for World War II, under the direct supervision of the Army Surgeon General. Its original location was at the Johns Hopkins School of Hygiene and Public Health. Its mission was to conduct occupational health surveys and investigations within the Department of Defense's (DOD's) industrial production base. It was staffed with three personnel and had a limited annual operating budget of three thousand dollars.

Most recently, it became internationally known as the U.S. Army Environmental Hygiene Agency (AEHA). Its mission expanded to support worldwide preventive medicine programs of the Army, DOD, and other Federal agencies as directed by the Army Medical Command or the Office of The Surgeon General, through consultations, support services, investigations, on-site visits, and training.

On 1 August 1994, AEHA was redesignated the U.S. Army Center for Health Promotion and Preventive Medicine with a provisional status and a commanding general officer. On 1 October 1995, the nonprovisional status was approved with a mission of providing preventive medicine and health promotion leadership, direction, and services for America's Army.

The organization's quest has always been one of excellence and the provision of quality service. Today, its goal is to be an established world-class center of excellence for achieving and maintaining a fit, healthy, and ready force. To achieve that end, the CHPPM holds firmly to its values which are steeped in rich military heritage:

★ *Integrity is the foundation*

★ *Excellence is the standard*

★ *Customer satisfaction is the focus*

★ *Its people are the most valued resource*

★ *Continuous quality improvement is the pathway*

This organization stands on the threshold of even greater challenges and responsibilities. It has been reorganized and reengineered to support the Army of the future. The CHPPM now has three direct support activities located in Fort Meade, Maryland; Fort McPherson, Georgia; and Fitzsimons Army Medical Center, Aurora, Colorado; to provide responsive regional health promotion and preventive medicine support across the U.S. There are also two CHPPM overseas commands in Landstuhl, Germany and Camp Zama, Japan who contribute to the success of CHPPM's increasing global mission. As CHPPM moves into the 21st Century, new programs relating to fitness, health promotion, wellness, and disease surveillance are being added. As always, CHPPM stands firm in its commitment to Army readiness. It is an organization proud of its fine history, yet equally excited about its challenging future.



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TRAINING MUNITIONS HEALTH RISK ASSESSMENT NO. 39-EJ-1485-00
RESIDENTIAL EXPOSURE FROM INHALATION OF AIR EMISSIONS
FROM THE LONG RIFLE .22 CALIBER BALL CARTRIDGE

EXECUTIVE SUMMARY

This assessment evaluated the potential for human health effects to offsite residents breathing air emissions following use of the Long Rifle .22 Caliber Ball Cartridge (.22 Caliber Ball) during training exercises.

To conduct this assessment, air emissions from the .22 Caliber Ball were collected in a test chamber at the U.S. Army Aberdeen Test Center, Maryland. The data collected from the Firing Point Emission Study provided the amount and types of substances released from the .22 Caliber Ball. This information was then used in an air dispersion model to determine ambient air concentrations at a location 100 meters (328 feet) downwind from a site where the .22 Caliber Ball may be used. Since the training facility in this assessment is hypothetical, the air model used assumptions that provided conservative estimates of air concentrations.

Modeled air concentrations were combined with exposure information (e.g., number of cartridges used per year) to estimate the amount of each substance the hypothetical offsite resident breathes. This estimate was then compared with the substance's health based screening level, which was obtained from agencies such as the U.S. Environmental Protection Agency, to determine if there is a potential for health effects from inhalation of these substances.

The health risk assessment included both long-term (30 years) and short-term (15-minute or 1-hour) exposures to modeled substance concentrations. Assessment results, generated using conservative methods, showed that the hypothetical offsite resident breathing air as close as 100 meters from the .22 Caliber Ball firing location is safe from these emissions. It should be noted that at most training installations, training areas are over 1,000 meters (over half a mile) away from populated areas.

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14. ABSTRACT

This assessment evaluated the potential for human health effects to offsite residents breathing air emissions following use of the Long Rifle .22 Caliber Ball cartridge. This document present the evaluation of the potential for adverse human health effects to the offsite residents breathing air emissions following the use of military firing ranges during training exercises. Study results showed no potential for health risks to the hypothetical resident from inhalation of air emissions from the .22 Caliber Cartridge. To conduct this study, air emissions from the .22 Caliber Ball Cartridge were collected in a test chamber (at Aberdeen Test Center, Aberdeen, MD). This information was then used in an air dispersion model to determine ambient air concentrations at a location downwind from the site where the item was activated. Modeled air concentrations were combined with exposure information to estimate the amount of substances the hypothetical resident breathes. This intake was combined with the substance's health information, to determine if there is a potential for health risks from inhalation of these substances. The health risk included both long-term and short term exposures to the modeled substance concentrations. Study results showed no potential for health risks from inhalation of air emissions from the .22 Caliber Ball Cartridge

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LIST OF ACRONYMS

| | |
|------------------|---|
| AEC | U.S. Army Environmental Center |
| AEGL | Acute Exposure Guideline Levels |
| AIHA | American Industrial Hygiene Association |
| ATC | U.S. Army Aberdeen Test Center |
| ATV | Acute Toxicity Value |
| DOE | U.S. Department of Energy |
| DODIC | Department of Defense Identification Code |
| EPA | U.S. Environmental Protection Agency |
| ERPG | Emergency Response Planning Guidelines |
| HBSL | Health-Based Screening Level |
| INPUFF | Integrated PUFF Model |
| NAAQS | National Ambient Air Quality Standards |
| NEW | Net Explosive Weight |
| OEL | Occupational Exposure Limit |
| PM ₁₀ | Particulate Matter under 10 microns in size |
| PRG | Preliminary Remediation Goals |
| RBC | Risk-Based Concentration |
| RfC | Reference Concentration |
| TEEL | Temporary Emergency Exposure Limits |
| TPH | Total Petroleum Hydrocarbons |
| TSP | Total Suspended Particulates |
| USACHPPM | U.S. Army Center for Health Promotion and Preventive Medicine |

TRAINING MUNITIONS HEALTH RISK ASSESSMENT NO. 39-EJ-1485-00
RESIDENTIAL EXPOSURE FROM INHALATION OF AIR EMISSIONS FROM THE
LONG RIFLE .22 CALIBER BALL CARTRIDGE

1. PURPOSE

This document presents the assessment of the potential for human health effects to offsite residents breathing air emissions following use of the Long Rifle .22 Caliber Ball Cartridge (.22 Caliber Ball) on firing ranges during training exercises.

2. AUTHORITY

Statement of Work, 30 November 2000, Training Munitions Inhalation Health Risk Evaluations.

3. REFERENCES

See Appendix A for a list of references.

4. BACKGROUND

4.1 CARTRIDGES AND THEIR USE

Cartridges are cases that contain a primer, propelling charge, and projectile. The primer is needed to activate the propelling charge, which provides the force to send the projectile to a target. Examples of projectiles include bullets, rockets, and missiles. Cartridges are also referred to as "rounds" and are fired from weapons such as pistols or rifles. The use of cartridges with weapons during training activities is important in preparing our soldiers for a variety of combat situations.

4.2 WHAT IS THE LONG RIFLE .22 CALIBER BALL CARTRIDGE?

The .22 Caliber Ball is a type of ball ammunition used for marksmanship practice and match use. The .22 Caliber Ball consists of a copper alloy cartridge case and a lead-antimony bullet. It also contains a propelling charge that consists mostly of nitrocellulose. Nitrocellulose is the primary ingredient in smokeless propellant (for both military and commercial use) and is also used in the production of lacquers and artificial leathers. Each .22 Caliber Ball cartridge is about as long as the width of a quarter (Reference 1).

4.3 ASSESSMENT SUMMARY

The .22 Caliber Ball was evaluated using an approach consisting of two main parts: air dispersion modeling and exposure assessment, which are briefly discussed in the paragraphs below. Sections 5 through 7 present a discussion of the methodology used for this assessment.

Emissions data used in the air dispersion modeling were obtained from the Firing Point Emission Study, conducted by the U.S. Army Aberdeen Test Center (ATC), at Aberdeen Proving Ground, Maryland (Reference 2). This study was funded by the U.S. Army Environmental Center (AEC) with the purpose of identifying and quantifying emissions from weapons firing. Data from this study were generated by firing munitions in a test chamber using weapons that are representative of those used by the U.S. Army during training. Emissions data for the .22 Caliber Ball were generated by firing it from either the Armalite® AR-7 Explorer Rifle or an M16 rifle fitted with a caliber .22 Ball, Long Rifle conversion kit.

The emissions data for the .22 Caliber Ball were used with an atmospheric dispersion model to estimate the average concentrations that may be experienced by an offsite resident. Since this assessment is designed to provide results that would be applicable to most Army training facilities, the training area used in this assessment was a hypothetical one. While most training areas are at least 1,000 meters away from populated areas, as a conservative distance, it was initially assumed that a person could reside 100 meters downwind from the firing point (location where the rifle is positioned). In addition, air-modeling parameters were selected to mimic worst-case conditions.

The exposure assessment included calculations of time-averaged concentrations for both long-term (chronic) and short-term (acute) exposures. For the purpose of this assessment, air concentrations were averaged over 30 years for chronic exposures and 1-hour or 15 minutes for acute exposures. Using a screening approach, a substance's estimated time-averaged air concentration was then compared to a chronic health-based screening level (HBSL) selected from sources established by the U.S. Environmental Protection Agency (EPA) and an acute toxicity value (ATV) selected from levels established by selected agencies depending on the exposure duration (i.e., 30 years versus 1-hour or 15 minutes). The terms HBSL and ATV are used for the purposes of this assessment. The comparison was made using the ratio of the HBSL or ATV to the estimated air concentration for each of the substances evaluated. If this ratio was less than one, no further evaluation was needed. This approach is conservative because the exposure assumptions used by the agencies, to establish HBSLs and ATVs, are likely to overestimate the exposures experienced by offsite residents living near firing ranges. If the chronic or acute averaged concentrations (C_{chronic} and C_{acute}) were greater than these screening levels, further analysis would be warranted to determine the potential for health effects. Note that concentrations greater than the screening levels do not indicate an onset of health effects, but rather the potential for such.

5. DATA COLLECTION AND AIR MODELING

5.1 EMISSION FACTORS

Emission factors, used to derive the air modeling emission rates used in this assessment, were generated from the Firing Point Emission Study conducted by the ATC (Reference 2). This study identified and quantified air emissions from the firing of

training munitions. The data included the net explosive weight (NEW), the substances sampled, and substance-specific emission factors. Emissions data from the Firing Point Emission Study are included in the first five columns of the table located in Appendix B.

5.2 BACKGROUND AND DESCRIPTION

Air dispersion models are available to mathematically simulate plume behavior and to estimate downwind concentrations of substances emitted from various sources. However, specific models are not available to determine the dispersion of emissions from munitions used during training. Estimating the magnitude and location of these concentrations depends on many factors including the amount and type of emissions, the behavior of the source, and meteorological conditions. Since a specific model is not available for modeling the use of munitions during training, the U.S. Army Center for Health Promotion and Preventive Medicine (USACHPPM) evaluated numerous air models to determine which would be suitable for use with munitions used during training. The USACHPPM recommended using the Integrated PUFF (INPUFF) model to estimate the dispersion of emissions from various munitions sources (Reference 3).

The INPUFF Model (Reference 4) was developed to simulate dispersion from instantaneous or semi-continuous point sources. This Gaussian-integrated puff model is capable of addressing a cloud type release over short periods of time, and computations can be performed for a single point source for multiple receptors. The algorithms used to calculate concentrations assume a vertically uniform wind direction (with no chemical reaction) to compute the contribution of each cloud at a receptor for each time step/interval.

5.3 MODEL ASSUMPTIONS

Some assumptions were made to best represent the firing of the .22 Caliber Ball cartridges. These assumptions were as follows:

- Typically, with conventional point sources (such as incinerators), the cloud rise and formation are determined by characterizing flue gas exit velocity, temperature, and stack diameter. However, the .22 Caliber Ball cartridges are used in conjunction with rifles. For unconventional sources with no real physical stack dimensions, such as rifles, the stack height and diameter were assumed to equal the height of the barrel and the bore diameter, respectively. No exit velocity was used with this source because the emission rates generated from the test data were obtained from sampling a stabilized cloud with no exit velocity. Table 1 includes the source parameters used to model the .22 Caliber Ball cartridges.

TABLE 1: SOURCE PARAMETERS

| Parameter | Model Input |
|--|--|
| Source/Stack Diameter | 0.009 meters |
| Source/Stack Height | 1 meter |
| Source Exit Temperature | 298.15 degrees Kelvin ($^{\circ}\text{K}$) (or 77 $^{\circ}\text{F}$) |
| Exit Velocity | 0 meters/second |
| Initial horizontal dispersion coefficient (σ_y) | 0.87 meters |
| Initial vertical dispersion coefficient (σ_z) | 1.07 meters |

- Initial cloud dimensions are preferred to model the air emissions from these types of releases. Typically, these dimensions are used to define the initial horizontal and vertical dispersion values (σ_y and σ_z) of the released cloud. However, this information was not measured during the studies at the ATC; therefore, the cloud dimensions were based on the test chamber dimensions and the volume of air sampled. By assuming an elliptical cloud with the prevailing wind direction being perpendicular to the rifle when fired, the test chamber's radius would be equal to the initial vertical dispersion (σ_z), and the initial horizontal dispersion (σ_y), would be equal to one half the length of the test chamber. The cloud exit temperature was assumed to be equal to the test chamber temperature.
- For the purposes of this assessment, a hypothetical offsite resident was assumed to be located 100 meters directly downwind from the source. The meander of the cloud is a major factor when estimating concentrations at given locations downwind from the source. Assuming that the resident is directly downwind from the source is the same as assuming that there is no cloud meander and the center of the cloud migrates directly over the hypothetical offsite resident. This assumption provides the most conservative modeled concentrations.
- Since this assessment does not look at a specific training site, generic, worst-case meteorological data were used. To determine the worst-case meteorological conditions that would result in the highest air emission concentrations, the modeling was performed using the EPA Risk Management Program Guidance (Reference 5). This guidance includes tables for estimating the footprint of chemical releases and is intended to inform emergency responders of potential accidental releases. The EPA has defined most default conditions for meteorological modeling parameters. Table 2 lists the meteorological parameters that were used in the air model.

TABLE 2: WORST-CASE METEOROLOGICAL PARAMETERS

| Parameter | Input Value |
|-----------------------|------------------------------------|
| Wind Speed | 1 meter/second |
| Atmospheric Stability | Category F |
| Wind Direction | 270° |
| Ambient Temperature | 293 degrees Kelvin (°K) (or 68 °F) |

5.4 GENERAL METHODOLOGY

The model was run for a total calculation time of 200 seconds to ensure that the total mass of the cloud had passed the hypothetical resident location. Concentrations were calculated every 2 seconds. The model results indicated that the initial cloud reached the hypothetical offsite resident within 80 seconds and dissipated below the lowest concentration the model calculated, which in this instance ($1 \times 10^{-11} \text{ g/m}^3$) occurred within 138 seconds. Table 3 contains the air model input parameters used in this assessment.

TABLE 3: AIR MODEL INPUT PARAMETERS

| Parameter | Input Value |
|---|-------------|
| Number of meteorological periods (NTIME) | 1 |
| Duration of each meteorological period (ITIME) | 200 seconds |
| Number of updates to the source (NSRCDS) | 100 |
| Duration/time step between each source update (ISUPDT) | 2 seconds |
| Total time modeled/Simulation Period (NTIME) (ITIME)= (NSRCDS) (ISUPDT) | 200 seconds |

5.5 USE OF MODEL OUTPUT

The concentrations provided by the INPUFF model were based on a unit emission rate (ER_{unit}) of 1 gram/second from an emission source, and did not represent any substance-specific concentrations from the use of any weapons system. This unit emission rate is typically used for ease of modeling purposes. The relationship between the emission rate and predicted concentration is linear. Therefore, the ratio of the predicted concentration to the unit emission rate was multiplied by each substance-specific emission rate to provide substance-specific concentrations.

5.6 DETERMINATION OF SUBSTANCE-SPECIFIC EMISSION RATES

The actual substance emission rate for one item (ER_1) for each substance was calculated using Equation 1. Example 1 contains a sample calculation using this equation.

$$ER_1 = \frac{EF \cdot CV}{t} \quad \text{Equation 1}$$

Where:

ER_1 = emission rate for one item (g/item)/sec
 EF = average adjusted emission factor (lb/item)
 CV = conversion factor (453.59 g/lb)
 t = release duration obtained from the INPUFF model (sec)

Example 1

Sample Calculation Using Equation 1:

$$ER_1 = \frac{(5.90E-08)(453.59)}{(2)}$$

$$= 1.339E-05 \text{ g/sec/item}$$

Calculation provided for benzene. Appendix B provides the average adjusted emission factor (EF) in lb/item.

Substance-specific ambient concentrations for one item (CONC) were calculated using Equation 2. A sample calculation using this equation is provided in Example 2. Appendix B contains the estimated air concentrations.

$$CONC = ER_1 \cdot \frac{UC}{ER_{unit}} \quad \text{Equation 2}$$

Where:

$CONC$ = substance concentration based on one item (g/m³)
 ER_1 = emission rate for one item (g/sec)
 ER_{unit} = unit emission rate as used in the model (g/sec)
 UC = concentration based on the unit emission rate (g/m³)

Example 2

Sample Calculation Using Equation 2:

$$CONC = (1.339E - 05) \frac{(2.061E - 04)}{(1)}$$

$$= 2.760 E-09 \text{ g/m}^3$$

Calculation provided for benzene.

6. RISK ASSESSMENT

6.1 EXPOSURE ASSUMPTIONS

Exposure assumptions were selected using a typical use scenario for the .22 Caliber Ball. The typical use scenario was provided by the AEC and is based on consultation with their senior training advisor (References 6, 7). The frequency of use for the .22 Caliber Ball was required to determine how much substance an offsite resident would be exposed to in the time period of interest (i.e., acute or chronic exposure). Table 4 summarizes the general use scenario for the .22 Caliber Ball.

TABLE 4: FREQUENCY OF USE FOR THE LONG RIFLE .22 CALIBER BALL CARTRIDGE

| Parameter | Value Used |
|---|------------|
| Number of cartridges used per year | 6,550 |
| Maximum number of cartridges used in 1-hour | 200 |

6.2 TIME-AVERAGING

For the chronic assessment, time-averaged concentrations were calculated by assuming that the hypothetical offsite resident would be exposed for 30 years. This is consistent with the exposure duration used by the EPA, which assumes that the resident spends 30 years at the same residence. By using the same exposure duration, the estimated time-averaged concentrations could be compared with their respective HBSLs, which are derived using standard EPA default assumptions.

Using the default residence time established by the EPA, the assumption was made that someone could be exposed to air emissions from 6,550 cartridges per year for 30 years. Table 5 lists the exposure parameters used to estimate concentrations for the chronic assessment. These parameters are based on the typical use scenario provided by AEC (Table 4) and the assumptions used in the air model run.

TABLE 5: EXPOSURE PARAMETERS USED TO DETERMINE TIME-AVERAGED CHRONIC AIR CONCENTRATIONS

| Exposure Parameter | Value Used |
|---|----------------------------------|
| Exposure Time (ET _{ctg}) | 3.333 min/cartridge ¹ |
| Exposure Frequency (EF _{ctg}) | 6,550 cartridges/year |
| Exposure Duration (ED) | 30 years ² |
| ¹ Based on the total model time of 200 seconds (3.33 minutes) used in the air model run. | |
| ² EPA default value. | |

Chronic averaged concentrations were calculated using Equation 4. Example 4 shows how this calculation was performed. Since benzene is classified as a carcinogen, as indicated in Appendix C, the averaging time (AT) is 70 years.

$$C_{chronic} = \frac{CONC \cdot 10^6 \cdot ET_{ctg} \cdot EF_{ctg} \cdot ED}{525,600 \cdot AT} \quad \text{Equation 4}$$

Where:

$C_{chronic}$ = average chronic concentration (µg/m³)
 $CONC$ = average modeled concentration for one cartridge (g/m³)
 10^6 = unit conversion (µg/g)
 ET_{ctg} = exposure time per cartridge (minutes/cartridge)
 EF_{ctg} = exposure frequency (cartridges/year)
 ED = exposure duration (years)
 $525,600$ = unit conversion (minutes/year)
 AT = averaging time (years)
 (Carcinogenic endpoint: AT = 70 years)
 Noncarcinogenic endpoint: AT = ED

Example 4
Sample Calculation Using Equation 4:

$$C_{chronic(benzene)} = \frac{(2.760E-09)(10^6)(3.333)(6,550)(30)}{(525,600)(70)}$$

$$= 4.91E-05 \mu\text{g}/\text{m}^3$$

The average modeled concentration for one cartridge (CONC) was obtained from Appendix B. The exposure parameters were obtained from Table 5.

Since many cartridges may be fired in a short period of time, acute exposures cannot be overlooked. Unlike the chronic assessment, only limited guidance for evaluating acute exposures is currently available. For the purpose of this assessment, acute exposure is defined as a 1-hour or 15-minute exposure. The 1-hour or 15-minute acute exposure averaging times allow for comparison with guidelines developed specifically for emergency planning purposes (see discussion on acute toxicity below).

The exposure frequency is based on the number of cartridges used per 1-hour or 15 minutes depending on the guideline used for comparison. This information is based on the use scenario provided by the AEC (Table 4). To estimate air concentrations for potential acute health impacts, it was conservatively assumed that 200 .22 Caliber Ball cartridges are fired in 1-hour. The average acute concentrations were computed using Equation 5. Example 5 contains a sample calculation using this equation. Benzene is used as the example substance.

$$C_{acute} = \frac{CONC \cdot 10^6 \cdot ET_{ctg} \cdot EF_{ctg}}{60} \quad \text{Equation 5}$$

Where:

- C_{acute} = average acute concentration ($\mu\text{g}/\text{m}^3$)
- CONC = average modeled concentration for one cartridge (g/m^3)
- 10^6 = unit conversion ($\mu\text{g}/\text{g}$)
- ET_{ctg} = exposure time per cartridge (minutes/cartridge)
- EF_{ctg} = exposure frequency (cartridges/hour)*
- 60 = unit conversion (minutes/hour)

* Based on 1-hour or 15 minute (0.25 hour) ATV

Example 5
Sample Calculation Using Equation 5:

$$C_{acute(benzene)} = \frac{(2.760E-09)(10^6)(3.333)(200)}{60}$$

$$= 3.07E-02 \mu\text{g}/\text{m}^3$$

The average modeled concentration for one cartridge (CONC) for benzene was obtained from Appendix B. See Appendix C to determine the ATV used.

6.3 TOXICITY ASSESSMENT

The potential for health effects was determined by comparing time-averaged air concentrations to health-based screening levels, which are developed from a substance's known toxicity. These toxicity values typically include different levels of safety factors depending on the level of confidence of the critical study. Appendix C contains a table of screening levels used for the chronic and acute assessments.

6.3.1 CHRONIC ASSESSMENT

The chronic assessment was conducted using a screening approach. Using this method, a substance's estimated time-averaged air concentration was compared to its HBSL. If this ratio was less than one, no further analysis was needed. This approach is conservative because the exposure assumptions used by the EPA, to establish HBSLs, assume that the resident is exposed for 350 days per year (assuming 2 weeks vacation per year). In contrast, exposure to air emissions from actual training activities at a firing range is intermittent and is not likely to occur on a daily basis year round.

A hierarchy of sources was developed for selection of the HBSLs to quantitatively evaluate as many of the identified substances as possible. The hierarchy of sources used was as follows:

- Clean Air Act, EPA National Ambient Air Quality Standards (NAAQS) (Reference 8)
- EPA Region 9 Preliminary Remediation Goals (PRGs) (Reference 9)
- EPA Region 3 Risk-Based Concentrations (RBCs) (Reference 10)

Some substances have neither PRGs nor RBCs because they have their own set of regulatory standards. Under the Clean Air Act, the EPA is required to establish NAAQS for several substances considered harmful to public health and the environment. Currently, NAAQS are available for seven substances. The NAAQS for the longer averaging time were used for the chronic assessment. Depending on the substance, this can range from an 8-hour average to an annual average. In addition, since the majority of the measured total suspended particulates (TSP) were PM₁₀ (particulate matter under 10 microns in size) (Reference 2), the NAAQS for PM₁₀ was used to evaluate the potential for health effects from exposure to TSP.

Next on the hierarchy, after the NAAQS, are the EPA Region 9 PRGs and the EPA Region 3 RBCs. The methodology used by EPA Region 9 to develop the PRGs generally results in lower values than the EPA Region 3 RBCs. However there were occasions when the RBCs were lower than the PRGs. To maintain a conservative approach for this assessment, the lower of the two values from these sources was selected as the HBSL for each substance evaluated. If only one value was available from these sources it was selected as the HBSL. To ensure that the most recent information was used, the Internet sites of both EPA Regions were checked. Appendix C presents the HBSLs used for this assessment.

Although the general approach used by both EPA Region 3 and Region 9 is the same, the exposure assumptions differ enough so that final recommended values can vary to a certain degree. In both methods, a substance's screening concentration was selected using the toxicity endpoint that derives a lower concentration. For example, if a substance has a known systemic toxicity and is a carcinogen, the screening concentration was calculated using both toxicity values. To maintain a conservative approach, EPA then selected the lower screening concentration as the recommended PRG or RBC.

Example 6 shows a sample calculation of how a substance's estimated chronic concentration was compared to its HBSL using benzene concentration.

Example 6

Sample Calculation Comparing a Substance's Estimated Chronic Concentration to Its HBSL:

$$\frac{C_{\text{chronic}(\text{benzene})}}{HBSL} = \frac{4.91E-05}{2.16E-01}$$
$$= 2.27E-04 < 1$$

In this case, the resulting ratio is less than one, indicating further evaluation is not necessary.

Many petroleum hydrocarbons were detected but do not have specific screening levels. Therefore, the approach recommended by the Total Petroleum Hydrocarbon Criteria Working Group (Reference 11) was adopted to evaluate petroleum hydrocarbon mixtures. Based on the working group's assessment of various hydrocarbons, it was recommended that mixtures be separated according to a substance's number of carbons and its chemical class (i.e., aliphatic or aromatic¹). Generally, as a substance's carbon number increases, its molecular weight increases, and it is, therefore, not a substance of concern via inhalation. The working group also concluded that aromatic hydrocarbons tend to be more toxic than aliphatic hydrocarbons (Reference 11). Table 6 tabulates the inhalation toxicity values used to evaluate exposure to petroleum mixtures. To be consistent with the methodology used in this assessment, the reference concentrations (RfCs) were converted to PRGs using Region 9 exposure assumptions. The resulting PRGs were used as the HBSLs for the petroleum hydrocarbons in this assessment. Appendix D presents these values.

¹ Aliphatic hydrocarbons are hydrocarbons in which the carbon atoms are joined by single covalent bonds consisting of two shared electrons (e.g., butane). Aromatic hydrocarbons have ring structures (e.g., benzene) (Reference 12).

TABLE 6: SUMMARY OF RfCs USED FOR PETROLEUM HYDROCARBONS¹

| Carbon Range | Aromatic Inhalation RfC (mg/m ³) | Aliphatic Inhalation RfC (mg/m ³) |
|--|---|--|
| C ₅ – C ₆ C _{>6} – C ₈ | | 18.4 |
| C _{>7} – C ₈ | 0.4 | |
| C _{>8} – C ₁₀ C _{>10} – C ₁₂ C _{>12} – C ₁₆ | 0.2 | 1.0 |
| C _{>16} – C ₂₁ C _{>21} – C ₃₅ | NA | NA |
| ¹ Reference 11 NA = not applicable for high molecular weight total petroleum hydrocarbons (C _{>16}) because substances in this carbon range are not volatile and therefore, inhalation is not a pathway of concern. | | |

6.3.2 ACUTE ASSESSMENT

An established method for assessing acute health effects is not currently available. In 1995 the EPA recognized the need for acute exposure guidelines for emergency response purposes and created the National Advisory Committee for Acute Exposure Guideline Levels (AEGLs) for Hazardous Substances. Currently, AEGLs are available for only a few substances.

To overcome the unavailability of acute toxicity data, several state regulatory agencies have suggested that guidelines developed for emergency purposes be used in the interim. Although suggestions have been made to use occupational exposure limits (OELs) by applying additional safety factors (References 13, 14), OELs were not used in this assessment because they introduce even more uncertainty than the use of emergency guidelines. The OELs are designed to protect the workplace environment and assume 8 hours a day, 5 days a week exposures. By definition, these exposures are more chronic than acute. In comparison, emergency planning guidelines are more appropriate because they are typically developed for exposures of 1-hour or less. In addition, safety factors are included as part of the guideline development so that the values would be protective of the general population.

For this study, the hierarchy of sources for ATV selection was as follows with each ATV defined below:

- EPA AEGL-1. "AEGL-1 is the airborne concentration of a substance above which it is predicted that the general population, including susceptible individuals, could experience notable discomfort, irritation, or certain asymptomatic, nonsensory effects. However, the effects are not disabling and are transient and reversible upon cessation of exposure."

- AIHA ERPG-1. "The maximum concentration in air below which it is believed nearly all individuals could be exposed for up to 1- hour without experiencing other than mild transient adverse health effects or perceiving a clearly defined objectionable odor."
- DOE TEEL-1. "The maximum concentration in air below which it is believed nearly all individuals could be exposed without experiencing other than mild transient adverse health effects or perceiving a clearly defined objectionable odor."

Emergency Response Planning Guidelines (ERPGs) published by the American Industrial Hygiene Association (AIHA) (Reference 15) and the Temporary Emergency Exposure Limits (TEELs) developed by the U.S. Department of Energy (DOE) (Reference 16) were used for this assessment, specifically the ERPG-1s and the TEEL-1s. Since TEEL-1s are intended for exposures up to 15-minutes, air concentrations compared to TEELs were averaged over a 15-minute period. Air concentrations compared to ERPGs and AEGLs were averaged over 1-hour, as these values are intended for 1-hour exposures.

AEGLs were used first when available since they are developed specifically for the purpose of acute exposure assessments. The ERPGs were selected next, prior to a substance's TEEL, because they are vigorously reviewed before they are published, whereas the TEELs are not.

Example 7 shows a sample calculation of how a substance's estimated acute concentration was compared to its ATV.

Example 7
Sample Calculation of Comparing a Substance's Estimated Acute Concentration to Its Acute Toxicity Value:

$$\frac{C_{acute(benzene)}}{ATV} = \frac{3.07E-02}{1.56E+05}$$
$$= 1.97E-07 < 1$$

In this example with benzene, the ratio is less than one, indicating that further analysis is not necessary.

7. RISK CHARACTERIZATION

As previously described, the exposure assessment included calculations of time-averaged concentrations for both long-term (chronic) and short-term (acute) exposures. Using a screening approach, a substance's estimated time-averaged air concentration was then compared to chronic HBSLs or ATVs. The comparison was made using the ratio of the HBSL or ATV to the estimated concentration. This approach is conservative because the exposure assumptions used by the EPA and other agencies, to establish HBSLs and ATVs, are likely to overestimate the exposures experienced by offsite residents living near firing ranges.

If this ratio was less than one, no further evaluation was needed. If the chronic or acute averaged concentrations (C_{chronic} and C_{acute}) were greater than the screening levels, resulting in a ratio greater than one, further evaluation would be warranted to determine the potential for health effects. Note that concentrations greater than the screening levels do not indicate an onset of health effects, but rather, the potential for such.

The chronic and acute assessments were conducted as outlined in Section 6.3. Appendix D presents results from the .22 Caliber Ball risk characterization.

7.1 CHRONIC HEALTH RISK

The outcome of the chronic assessment indicated that no chronic health effects are expected from breathing the air emissions from the .22 Caliber Ball. Since the ratios for all substances were below one, further evaluation was not needed.

7.2 ACUTE HEALTH RISK

For the acute assessment, all ratios were below one, indicating that no acute health effects are expected from breathing the air emissions from the .22 Caliber Ball. The ratios for all substances were less than one, indicating further evaluation was not necessary.

7.3 FACT SHEET

Appendix E includes a copy of the fact sheet submitted to the AEC. The fact sheet uses the results from this assessment to communicate information related to inhalation of .22 Caliber Ball air emissions.

8. UNCERTAINTY DISCUSSION

The limitations inherent in modeling and the added conservatism of the assessment contribute to the uncertainty of the assessment results. The risk assessment methodology typically includes safety factors that are embedded in the toxicity data to ensure adequate protection of the general population, particularly, susceptible individuals such as the sick, elderly, and children. Table 7 identifies areas of uncertainty associated with this assessment.

TABLE 7: TYPES OF UNCERTAINTY

| Issue | Uncertainty | Direction of Effect |
|---|--|---------------------|
| Modeling | | |
| Modeled versus real-time sampling | The air concentrations in this assessment were modeled. Actual air concentrations taken from the field may be higher or lower. | Varies |
| Frequency of use for the .22 Caliber Ball | Actual frequency of use for these munitions during training exercises may be different from those stated in this report. | Varies |
| Hypothetical offsite resident assumed to be located directly downwind | Unless the area around the training facility is populated, the chances that a person living directly downwind is low. | Overestimates |
| Use of worst-case meteorological conditions | To ensure that this assessment is applicable to most training areas, worst-case meteorological conditions were used in the air model. | Overestimates |
| Exposure Assessment | | |
| Estimating time-averaged concentrations | Actual exposure from the .22 Caliber Ball is intermittent. If one were to plot a person's exposure profile, the plot would consist of a series of spikes. Since current risk assessment methodology does not allow the evaluation of the potential for health risks as a function of time, a single concentration, averaged over the exposure duration was used. In this assessment, the exposure durations used were 30 years and 1-hour or 15 minutes. | Varies |
| Comparing estimated concentration to established screening levels | The Region 3 and Region 9 HBSLs were developed using different exposure assumptions than those in this assessment, resulting in more conservative screening levels. | Overestimates |
| Comparing estimated concentrations to established screening levels | Comparison to screening levels does not account for possible cumulative effects of exposure to more than one substance. | Underestimates |

TABLE 7: TYPES OF UNCERTAINTY

| Issue | Uncertainty | Direction of Effect |
|---|--|----------------------------|
| Screening assessment versus calculating an average daily intake | Calculating an average daily intake allows the use of scenario-specific assumptions. However, unless the ratio of concentration to screening level approaches one, a screening assessment is useful as a first-cut evaluation. | Varies |
| Exposure to other munitions | Other munitions are typically used during the same training exercise. These items may contain similar or different substances from those detected in the .22 Caliber Ball. | Underestimates |
| Toxicity Assessment | | |
| Lack of toxicity data | Some substances were not quantitatively evaluated because they have no known toxicity data. | Underestimates |
| Modifying and uncertainty factors for toxicity data | Modifying factors and uncertainty factors of varying degree are typically applied to toxicological values. These factors are used to conservatively account for extrapolating from animal studies for human health evaluation, and to conservatively account for variation in human populations. | Overestimates |

9. CONCLUSION

Using conservative assumptions, the assessment indicated that offsite residents who live as close as 100 meters directly downwind from the firing location are safe from breathing air emissions from the .22 Caliber Ball. It is believed that the assumptions contained in this analysis are conservative enough to be protective of all the population including the sick, elderly, and children.

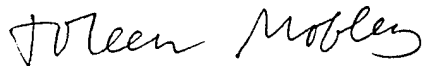
10. RECOMMENDATIONS

The results from this assessment are intended for a hypothetical training facility, and actual results may vary depending on site-specific conditions. This assessment used conservative assumptions (e.g., worst-case meteorological conditions, receptor located directly downwind, etc.) and it is believed that most site-specific analyses would result in even lower concentrations. Therefore, the results from this assessment should be applicable to most training facilities, unless site-specific conditions vary significantly.

10. POINT OF CONTACT

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APPENDIX A
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APPENDIX B

AIR DISPERSION MODELING OUTPUT DATA

Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Cartridge, Caliber .22 Ball, Long Rifle | | | | | | | No. of rounds (l) | | 1 round | |
|--|--|--|--|--|--|--|---|---|-----------------------------------|--|
| Number of items: Trial #1B => 71 Trial #2B => 71 | | | | | | | release duration (t): | | 2 seconds | |
| Net Explosive Weight- N.E.W. per item (lbs.) => 3.57E-04 | | | | | | | Unit Concentration (UC): | | 2.061E-04 g/m ³ /(g/s) | |
| ATC Firing Test Results ¹ | | | | | | | | | | |
| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER ₁ | | |
| | | | | | | | | | | |
| Permanent Gases | | | | | | | | | | |
| Ammonia (NH ₃) | 7.00E+00 | 7.00E+00 | NA | ND | ND | ND | ND | ND | | |
| Carbon Dioxide (CO ₂) | 3.62E+02 | 3.49E+02 | NA | 7.38E-05 | 2.07E-01 | 3.347E-02 | 3.449E-06 | 1.674E-02 | | |
| Carbon Monoxide (CO) | 3.80E+02 | 3.81E+02 | NA | 7.89E-05 | 2.21E-01 | 3.579E-02 | 3.688E-06 | 1.789E-02 | | |
| Oxides of Nitrogen (NO _x) | 1.54E+01 | 1.56E+01 | NA | 3.22E-06 | 9.01E-03 | 1.459E-03 | 1.504E-07 | 7.296E-04 | | |
| Sulfur Dioxide (SO ₂) | 5.24E-01 | 5.24E-01 | NA | ND | ND | ND | ND | ND | | |
| Acid Gases | | | | | | | | | | |
| Hydrogen Fluoride | 2.30E-01 | 2.30E-01 | 2.50E-01 | ND | ND | ND | ND | ND | | |
| Hydrogen Chloride | 2.30E-01 | 2.20E-01 | 2.40E-01 | ND | ND | ND | ND | ND | | |
| Hydrogen Bromide | 2.20E-01 | 2.20E-01 | 2.40E-01 | ND | ND | ND | ND | ND | | |
| Nitric Acid | 2.30E-01 | 3.70E-01 | 2.40E-01 | 8.71E-08 | 2.44E-04 | 3.949E-05 | 4.069E-09 | 1.974E-05 | | |
| Phosphoric Acid | 2.30E-01 | 2.20E-01 | 2.40E-01 | ND | ND | ND | ND | ND | | |
| Sulfuric Acid | 2.30E-01 | 2.20E-01 | 2.40E-01 | ND | ND | ND | ND | ND | | |
| Cyanide | | | | | | | | | | |
| Particulate Cyanide | 1.30E-02 | 1.30E-02 | 1.40E-02 | ND | ND | ND | ND | ND | | |
| Hydrogen Cyanide | 2.67E-01 | 4.11E-01 | 1.40E-02 | 8.15E-08 | 2.28E-04 | 3.696E-05 | 3.808E-09 | 1.848E-05 | | |
| Particulate | | | | | | | | | | |
| Total Suspended Particulate | 1.27E+01 | 1.41E+01 | NA | 3.25E-06 | 9.08E-03 | 1.472E-03 | 1.517E-07 | 7.360E-04 | | |
| Particulate Matter <10 microns | 1.30E+01 | 1.43E+01 | NA | 3.30E-06 | 9.22E-03 | 1.495E-03 | 1.540E-07 | 7.473E-04 | | |
| Particulate Matter <2.5 microns | 9.92E+00 | 1.10E+01 | NA | 2.52E-06 | 7.06E-03 | 1.143E-03 | 1.178E-07 | 5.716E-04 | | |
| Metals | | | | | | | | | | |
| Aluminum | 9.04E-02 | 1.74E-01 | 1.83E-01 | ND | ND | ND | ND | ND | | |
| Antimony | 3.50E-02 | 3.70E-02 | 1.17E-02 | 8.69E-09 | 2.43E-05 | 3.943E-06 | 4.063E-10 | 1.971E-06 | | |
| Arsenic | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND | | |
| Barium | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND | | |
| Beryllium | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND | | |

Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M _e | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER _i |
|-----------------------------------|--|--|--|--|--|---|---|---|
| Cadmium | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Calcium | 1.54E-01 | 2.71E-01 | 2.68E-01 | ND | ND | ND | ND | ND |
| Chromium | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Cobalt | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Copper | 2.61E-02 | 2.87E-02 | 1.17E-02 | 6.62E-09 | 1.85E-05 | 3.005E-06 | 3.096E-10 | 1.502E-06 |
| Lead | 7.20E+00 | 7.90E+00 | 1.17E-02 | 1.82E-06 | 5.10E-03 | 8.269E-04 | 8.521E-08 | 4.134E-04 |
| Magnesium | 3.21E-02 | 5.84E-02 | 1.63E-01 | ND | ND | ND | ND | ND |
| Manganese | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Nickel | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Selenium | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Silver | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Thallium | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Vanadium | 1.15E-02 | 1.17E-02 | 1.17E-02 | ND | ND | ND | ND | ND |
| Zinc | 6.27E-02 | 6.39E-02 | 1.17E-02 | 1.53E-08 | 4.28E-05 | 6.939E-06 | 7.150E-10 | 3.469E-06 |
| TO-11 Carbonyls | | | | | | | | |
| Formaldehyde | 5.30E-01 | 5.70E-01 | 2.20E-01 | 8.57E-08 | 2.40E-04 | 3.888E-05 | 4.007E-09 | 1.944E-05 |
| Acetaldehyde | 3.60E-01 | 3.50E-01 | 2.80E-01 | 2.59E-08 | 7.25E-05 | 1.176E-05 | 1.211E-09 | 5.878E-06 |
| Acetone | 1.00E+00 | 1.00E+00 | 1.00E+00 | ND | ND | ND | ND | ND |
| Acrolein | 2.00E-01 | 2.00E-01 | 2.00E-01 | ND | ND | ND | ND | ND |
| Propionaldehyde | 2.00E-01 | 2.00E-01 | 2.00E-01 | ND | ND | ND | ND | ND |
| Crotonaldehyde | 2.00E-01 | 2.00E-01 | 2.00E-01 | ND | ND | ND | ND | ND |
| Butyraldehyde | 1.59E+00 | 1.43E+00 | 1.55E+00 | 3.38E-08 | 9.46E-05 | 1.532E-05 | 1.579E-09 | 7.661E-06 |
| Benzaldehyde | 2.00E-01 | 2.00E-01 | 2.00E-01 | ND | ND | ND | ND | ND |
| Isovaleraldehyde | 2.00E-01 | 2.00E-01 | 2.00E-01 | ND | ND | ND | ND | ND |
| Valeraldehyde | 2.00E-01 | 2.00E-01 | 2.00E-01 | ND | ND | ND | ND | ND |
| o,m,p-Tolualdehyde | 6.00E-01 | 6.00E-01 | 6.00E-01 | ND | ND | ND | ND | ND |
| Hexaldehyde | 2.00E-01 | 2.00E-01 | 2.00E-01 | ND | ND | ND | ND | ND |
| 2,5-Dimethylbenzaldehyde | 2.00E-01 | 2.00E-01 | 2.00E-01 | ND | ND | ND | ND | ND |
| TO-14 VOCs (extended list) | | | | | | | | |
| Propene | 2.93E-01 | 3.10E-01 | 8.61E-04 | 6.89E-08 | 1.93E-04 | 3.123E-05 | 3.218E-09 | 1.562E-05 |
| Dichlorodifluoromethane | 2.97E-03 | 2.97E-03 | 3.46E-03 | ND | ND | ND | ND | ND |
| Chlorodifluoromethane | 3.54E-03 | 3.54E-03 | 3.54E-03 | ND | ND | ND | ND | ND |
| Freon 114 | 6.99E-03 | 6.99E-03 | 6.99E-03 | ND | ND | ND | ND | ND |

Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER ₁ |
|--------------------------|--|--|--|--|--|--|---|---|
| Chloromethane | 1.24E-03 | 1.45E-03 | 1.45E-03 | 7.71E-12 | 2.16E-08 | 3.496E-09 | 3.603E-13 | 1.748E-09 |
| Vinyl Chloride | 2.56E-03 | 2.56E-03 | 2.56E-03 | ND | ND | ND | ND | ND |
| 1,3-Butadiene | 3.98E-02 | 6.42E-02 | 2.21E-03 | 1.19E-08 | 3.34E-05 | 5.415E-06 | 5.580E-10 | 2.707E-06 |
| Bromomethane | 3.88E-03 | 3.88E-03 | 3.88E-03 | ND | ND | ND | ND | ND |
| Chloroethane | 2.64E-03 | 2.64E-03 | 2.64E-03 | ND | ND | ND | ND | ND |
| Dichlorofluoromethane | 4.21E-03 | 4.21E-03 | 4.21E-03 | ND | ND | ND | ND | ND |
| Trichlorofluoromethane | 1.69E-03 | 1.69E-03 | 1.69E-03 | 3.63E-11 | 1.02E-07 | 1.648E-08 | 1.698E-12 | 8.240E-09 |
| Pentane | 1.48E-03 | 1.19E-03 | 1.18E-03 | 5.90E-11 | 1.65E-07 | 2.674E-08 | 2.756E-12 | 1.337E-08 |
| Acrolein | 9.86E-02 | 1.10E-01 | 2.29E-03 | 2.39E-08 | 6.70E-05 | 1.085E-05 | 1.118E-09 | 5.424E-06 |
| 1,1-Dichloroethene | 4.05E-03 | 4.05E-03 | 4.05E-03 | ND | ND | ND | ND | ND |
| Freon 113 | 7.68E-03 | 7.68E-03 | 7.68E-03 | ND | ND | ND | ND | ND |
| Acetone | 3.09E-02 | 3.09E-02 | 3.09E-02 | 6.66E-10 | 1.86E-06 | 3.019E-07 | 3.111E-11 | 1.510E-07 |
| Methyl Iodide | 5.81E-03 | 5.81E-03 | 5.81E-03 | ND | ND | ND | ND | ND |
| Carbon Disulfide | 3.11E-03 | 3.11E-03 | 3.11E-03 | ND | ND | ND | ND | ND |
| Acetonitrile | 1.63E-01 | 9.23E-02 | 9.23E-02 | 1.00E-08 | 2.80E-05 | 4.535E-06 | 4.673E-10 | 2.267E-06 |
| 3-Chloropropene | 3.13E-03 | 3.13E-03 | 3.13E-03 | ND | ND | ND | ND | ND |
| Methylene Chloride | 1.35E+00 | 6.60E-01 | 2.43E-01 | 1.80E-07 | 5.03E-04 | 8.148E-05 | 8.397E-09 | 4.074E-05 |
| tert-Butyl Alcohol | 3.03E-03 | 3.03E-03 | 2.12E-03 | ND | ND | ND | ND | ND |
| Acrylonitrile | 3.04E-02 | 3.04E-02 | 2.17E-03 | 6.96E-09 | 1.95E-05 | 3.158E-06 | 3.254E-10 | 1.579E-06 |
| trans-1,2-Dichloroethene | 3.96E-03 | 3.96E-03 | 3.96E-03 | ND | ND | ND | ND | ND |
| Methyl t-Butyl Ether | 1.44E-03 | 3.61E-03 | 1.44E-03 | 2.82E-11 | 7.89E-08 | 1.278E-08 | 1.317E-12 | 6.388E-09 |
| Hexane | 2.47E-02 | 1.06E-02 | 3.17E-02 | ND | ND | ND | ND | ND |
| 1,1-Dichloroethane | 3.97E-03 | 3.97E-03 | 3.97E-03 | ND | ND | ND | ND | ND |
| Vinyl Acetate | 3.52E-03 | 3.52E-03 | 3.52E-03 | ND | ND | ND | ND | ND |
| cis-1,2-Dichloroethene | 3.96E-03 | 3.96E-03 | 3.96E-03 | ND | ND | ND | ND | ND |
| 2-Butanone | 2.95E-03 | 2.65E-03 | 2.95E-03 | 6.42E-10 | 1.80E-06 | 2.911E-07 | 2.999E-11 | 1.455E-07 |
| Ethyl Acetate | 3.60E-03 | 3.60E-03 | 3.60E-03 | ND | ND | ND | ND | ND |
| Methyl Acrylate | 3.52E-03 | 3.52E-03 | 3.52E-03 | ND | ND | ND | ND | ND |
| Chloroform | 4.88E-03 | 4.88E-03 | 4.88E-03 | ND | ND | ND | ND | ND |
| 1,1,1-Trichloroethane | 1.09E-02 | 1.64E-02 | 1.09E-02 | 8.66E-10 | 2.42E-06 | 3.928E-07 | 4.047E-11 | 1.964E-07 |
| Carbon Tetrachloride | 6.29E-03 | 6.29E-03 | 6.29E-03 | ND | ND | ND | ND | ND |
| 1,2-Dichloroethane | 4.05E-03 | 4.05E-03 | 4.05E-03 | 9.27E-10 | 2.60E-06 | 4.207E-07 | 4.335E-11 | 2.103E-07 |
| Benzene | 2.56E-01 | 2.62E-01 | 1.28E-03 | 5.90E-08 | 1.65E-04 | 2.678E-05 | 2.760E-09 | 1.339E-05 |

22cal_ball

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Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER ₁ |
|---------------------------|--|--|--|--|--|--|---|---|
| Isocutane | 4.67E-03 | 4.67E-03 | 4.67E-03 | ND | ND | ND | ND | ND |
| Heptane | 4.10E-03 | 4.10E-03 | 1.64E-02 | ND | ND | ND | ND | ND |
| Trichloroethane | 4.88E-03 | 4.88E-03 | 4.88E-03 | ND | ND | ND | ND | ND |
| Ethyl Acrylate | 4.09E-03 | 4.09E-03 | 4.09E-03 | ND | ND | ND | ND | ND |
| 1,2-Dichloropropane | 4.62E-03 | 4.62E-03 | 4.62E-03 | ND | ND | ND | ND | ND |
| Methyl Methacrylate | 4.09E-03 | 4.09E-03 | 4.09E-03 | ND | ND | ND | ND | ND |
| Dibromomethane | 7.11E-03 | 7.11E-03 | 7.11E-03 | ND | ND | ND | ND | ND |
| 1,4-Dioxane | 3.60E-03 | 3.60E-03 | 3.60E-03 | ND | ND | ND | ND | ND |
| Bromodichloromethane | 6.70E-03 | 6.70E-03 | 6.70E-03 | ND | ND | ND | ND | ND |
| cis-1,3-Dichloropropene | 4.54E-03 | 4.54E-03 | 4.54E-03 | ND | ND | ND | ND | ND |
| 4-Methyl-2-Pentanone | 4.10E-03 | 4.10E-03 | 4.10E-03 | ND | ND | ND | ND | ND |
| Toluene | 3.02E-02 | 2.64E-02 | 2.26E-03 | 6.00E-09 | 1.68E-05 | 2.723E-06 | 2.806E-10 | 1.362E-06 |
| Octane | 4.67E-03 | 4.67E-03 | 4.67E-03 | ND | ND | ND | ND | ND |
| trans-1,3-Dichloropropene | 4.54E-03 | 4.54E-03 | 4.54E-03 | ND | ND | ND | ND | ND |
| Ethyl Methacrylate | 4.67E-03 | 4.67E-03 | 4.67E-03 | ND | ND | ND | ND | ND |
| 1,1,2-Trichloroethane | 5.46E-03 | 5.46E-03 | 5.46E-03 | ND | ND | ND | ND | ND |
| Tetrachloroethane | 6.78E-03 | 6.78E-03 | 6.78E-03 | ND | ND | ND | ND | ND |
| 2-Hexanone | 4.10E-03 | 4.10E-03 | 4.10E-03 | ND | ND | ND | ND | ND |
| Dibromochloromethane | 8.52E-03 | 8.52E-03 | 8.52E-03 | ND | ND | ND | ND | ND |
| 1,2-Dibromoethane | 7.68E-03 | 7.68E-03 | 7.68E-03 | ND | ND | ND | ND | ND |
| Chlorobenzene | 4.60E-03 | 4.60E-03 | 4.60E-03 | ND | ND | ND | ND | ND |
| 1,1,1,2-Tetrachloroethane | 6.87E-03 | 6.87E-03 | 6.87E-03 | ND | ND | ND | ND | ND |
| Ethylbenzene | 8.68E-03 | 8.68E-03 | 8.68E-03 | ND | ND | ND | ND | ND |
| m/p-Xylene | 4.34E-02 | 3.04E-02 | 4.34E-02 | ND | ND | ND | ND | ND |
| o-Xylene | 4.78E-02 | 3.04E-02 | 5.65E-02 | ND | ND | ND | ND | ND |
| Styrene | 1.28E-02 | 1.28E-02 | 4.26E-03 | 2.93E-09 | 8.20E-06 | 1.328E-06 | 1.369E-10 | 6.642E-07 |
| Bromoform | 1.03E-02 | 1.03E-02 | 1.03E-02 | ND | ND | ND | ND | ND |
| Cumene | 1.97E-03 | 1.47E-03 | 1.97E-03 | ND | ND | ND | ND | ND |
| 1,1,2,2-Tetrachloroethane | 6.87E-03 | 6.87E-03 | 6.87E-03 | ND | ND | ND | ND | ND |
| 1,2,3-Trichloropropane | 6.03E-03 | 6.03E-03 | 6.03E-03 | ND | ND | ND | ND | ND |
| Bromobenzene | 6.42E-03 | 6.42E-03 | 6.42E-03 | ND | ND | ND | ND | ND |
| 4-Ethyltoluene | 3.93E-03 | 2.95E-03 | 3.93E-03 | ND | ND | ND | ND | ND |
| 1,3,5-Trimethylbenzene | 4.92E-03 | 3.93E-03 | 4.92E-03 | ND | ND | ND | ND | ND |

Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER ₁ |
|--------------------------|--|--|--|--|--|--|---|---|
| Alpha Methyl Styrene | 4.83E-03 | 4.83E-03 | 4.83E-03 | ND | ND | ND | ND | ND |
| 1,2,4-Trimethylbenzene | 1.47E-02 | 9.83E-03 | 1.47E-02 | ND | ND | ND | ND | ND |
| 1,3-Dichlorobenzene | 6.01E-03 | 6.01E-03 | 6.01E-03 | ND | ND | ND | ND | ND |
| 1,4-Dichlorobenzene | 6.01E-03 | 6.01E-03 | 6.01E-03 | ND | ND | ND | ND | ND |
| Benzyl Chloride | 5.18E-03 | 5.18E-03 | 5.18E-03 | ND | ND | ND | ND | ND |
| 1,2-Dichlorobenzene | 6.01E-03 | 6.01E-03 | 6.01E-03 | ND | ND | ND | ND | ND |
| Hexachlorethane | 9.68E-03 | 9.68E-03 | 9.68E-03 | ND | ND | ND | ND | ND |
| 1,2,4-Trichlorobenzene | 7.42E-03 | 7.42E-03 | 7.42E-03 | ND | ND | ND | ND | ND |
| Hexachlorobutadiene | 1.07E-02 | 1.07E-02 | 1.07E-02 | ND | ND | ND | ND | ND |
| Hydrocarbons | | | | | | | | |
| Methane | 3.35E+00 | 3.61E+00 | 1.34E+00 | 5.53E-07 | 1.55E-03 | 2.509E-04 | 2.586E-08 | 1.255E-04 |
| Ethylene | 1.46E+00 | 1.72E+00 | 2.87E-02 | 3.83E-07 | 1.07E-03 | 1.738E-04 | 1.791E-08 | 8.691E-05 |
| Acetylene | 4.54E-01 | 5.24E-01 | 2.66E-02 | 1.18E-07 | 3.30E-04 | 5.349E-05 | 5.512E-09 | 2.675E-05 |
| Ethane | 1.22E-01 | 1.46E-01 | 3.07E-02 | 3.23E-08 | 9.05E-05 | 1.466E-05 | 1.511E-09 | 7.331E-06 |
| Propylene | 2.34E-01 | 2.89E-01 | 4.30E-02 | 6.31E-08 | 1.77E-04 | 2.860E-05 | 2.947E-09 | 1.430E-05 |
| Propane | 4.51E-02 | 4.51E-02 | 4.51E-02 | ND | ND | ND | ND | ND |
| Propyne | 4.00E-02 | 4.00E-02 | 4.00E-02 | ND | ND | ND | ND | ND |
| Isobutane | 5.94E-02 | 5.94E-02 | 5.94E-02 | ND | ND | ND | ND | ND |
| 1-Butene/Isobutylene | 1.08E-01 | 1.08E-01 | 1.08E-01 | ND | ND | ND | ND | ND |
| 1,3-Butadiene/butane | 1.65E-01 | 1.65E-01 | 1.65E-01 | ND | ND | ND | ND | ND |
| cis-butene | 5.74E-02 | 5.74E-02 | 5.74E-02 | ND | ND | ND | ND | ND |
| 1-Butyne/trans-butene | 1.06E-01 | 1.06E-01 | 1.06E-01 | ND | ND | ND | ND | ND |
| 2-Butyne | 5.53E-02 | 5.53E-02 | 5.53E-02 | ND | ND | ND | ND | ND |
| n-Pentane | 7.38E-02 | 7.38E-02 | 7.38E-02 | ND | ND | ND | ND | ND |
| n-Hexane | 8.81E-02 | 8.81E-02 | 8.81E-02 | ND | ND | ND | ND | ND |
| SVOCs (8270 List) | | | | | | | | |
| N-nitrosodimethylamine | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Bis(2-chloroethyl)ether | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Phenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2-chlorophenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 1,3-dichlorobenzene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 1,4-dichlorobenzene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 1,2-dichlorobenzene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |

Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER ₁ |
|-----------------------------|--|--|--|--|--|--|---|---|
| Benzyl alcohol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Bis(2-chloroisopropyl)ether | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2-methylphenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Hexachloroethane | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| N-nitroso-di-n-propylamine | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 4-methylphenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Nitrobenzene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Isophorone | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2-nitrophenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2,4-dimethylphenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Bis(2-chloroethoxy)methane | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2,4-dichlorophenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 1,2,4-trichlorobenzene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Naphthalene | 1.53E-02 | 1.74E-02 | 1.80E-02 | 3.75E-09 | 1.05E-05 | 1.703E-06 | 1.755E-10 | 8.514E-07 |
| 4-chloroaniline | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Hexachlorobutadiene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 4-chloro-3-methylphenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2-methylnaphthalene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Hexachlorocyclopentadiene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2,4,6-trichlorophenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2,4,5-trichlorophenol | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2-chloronaphthalene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2-nitroaniline | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Acenaphthylene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Dimethylphthalate | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2,6-dinitrotoluene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Acenaphthene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 3-nitroaniline | 3.56E-02 | 3.59E-02 | 3.59E-02 | ND | ND | ND | ND | ND |
| 2,4-dinitrophenol | 3.56E-02 | 3.59E-02 | 3.59E-02 | ND | ND | ND | ND | ND |
| Dibenzofuran | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 2,4-dinitrotoluene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 4-nitrophenol | 3.56E-02 | 3.59E-02 | 3.59E-02 | ND | ND | ND | ND | ND |
| Fluorene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |

Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER ₁ |
|----------------------------|--|--|--|--|--|--|---|---|
| 4-chlorophenyl-phenylether | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Diethylphthalate | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 4-nitroaniline | 3.56E-02 | 3.59E-02 | 3.59E-02 | ND | ND | ND | ND | ND |
| 4,6-dinitro-2-methylphenol | 3.56E-02 | 3.59E-02 | 3.59E-02 | ND | ND | ND | ND | ND |
| N-nitrosodiphenylamine(1) | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| 4-bromophenyl-phenylether | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Hexachlorobenzene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Pentachlorophenol | 3.56E-02 | 3.59E-02 | 3.59E-02 | ND | ND | ND | ND | ND |
| Phenanthrene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Anthracene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Di-n-butylphthalate | 1.78E-02 | 1.29E-02 | 1.80E-02 | 2.99E-09 | 8.37E-06 | 1.356E-06 | 1.398E-10 | 6.782E-07 |
| Fluoranthene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Pyrene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Butylbenzylphthalate | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Benzo(a)anthracene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Chrysene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Bis(2-ethylhexyl)phthalate | 7.48E-02 | 7.19E-02 | 5.57E-02 | 5.24E-09 | 1.47E-05 | 2.378E-06 | 2.450E-10 | 1.189E-06 |
| Di-n-octylphthalate | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Benzo(b)fluoranthene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Benzo(k)fluoranthene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Benzo(a)pyrene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Indeno(1,2,3-cd)pyrene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Dibenz(a,h)anthracene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| Benzo(g,h,i)perylene | 1.78E-02 | 1.80E-02 | 1.80E-02 | ND | ND | ND | ND | ND |
| TO-13 PAHs | | | | | | | | |
| Naphthalene | 1.35E-02 | 1.56E-02 | 4.31E-04 | 3.25E-09 | 9.11E-06 | 1.476E-06 | 1.521E-10 | 7.381E-07 |
| Acenaphthylene | 1.60E-03 | 1.80E-03 | 3.77E-05 | 3.82E-10 | 1.07E-06 | 1.732E-07 | 1.785E-11 | 8.661E-08 |
| Acenaphthene | 1.07E-04 | 1.11E-04 | 3.23E-05 | 1.83E-11 | 5.12E-08 | 8.301E-09 | 8.554E-13 | 4.150E-09 |
| Fluorene | 2.32E-04 | 2.52E-04 | 5.03E-05 | 4.49E-11 | 1.26E-07 | 2.038E-08 | 2.100E-12 | 1.019E-08 |
| Phenanthrene | 3.21E-04 | 4.13E-04 | 1.47E-04 | 5.36E-11 | 1.50E-07 | 2.431E-08 | 2.505E-12 | 1.215E-08 |
| Anthracene | 3.56E-05 | 4.67E-05 | 1.80E-05 | 9.45E-12 | 2.64E-08 | 4.284E-09 | 4.415E-13 | 2.142E-09 |
| Fluoranthene | 7.30E-05 | 9.88E-05 | 3.77E-05 | 1.19E-11 | 3.33E-08 | 5.390E-09 | 5.555E-13 | 2.695E-09 |
| Pyrene | 1.16E-04 | 1.31E-04 | 2.69E-05 | 2.27E-11 | 6.36E-08 | 1.030E-08 | 1.062E-12 | 5.152E-09 |

Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER ₁ |
|---------------------------|--|--|--|--|--|--|---|---|
| Benzo(a)anthracene | 3.03E-05 | 3.77E-05 | 1.80E-05 | 7.80E-12 | 2.18E-08 | 3.538E-09 | 3.646E-13 | 1.769E-09 |
| Chrysene | 3.21E-05 | 4.67E-05 | 1.80E-05 | 9.04E-12 | 2.53E-08 | 4.101E-09 | 4.226E-13 | 2.050E-09 |
| Benzo(b)fluoranthene | 6.95E-05 | 8.26E-05 | 1.80E-05 | 1.74E-11 | 4.88E-08 | 7.912E-09 | 8.153E-13 | 3.956E-09 |
| Benzo(k)fluoranthene | 5.70E-05 | 6.83E-05 | 1.80E-05 | 1.44E-11 | 4.02E-08 | 6.516E-09 | 6.714E-13 | 3.258E-09 |
| Benzo(e)pyrene | 1.37E-04 | 1.42E-04 | 1.80E-05 | 3.20E-11 | 8.95E-08 | 1.451E-08 | 1.495E-12 | 7.253E-09 |
| Benzo(a)pyrene | 1.28E-04 | 1.44E-04 | 1.80E-05 | 3.12E-11 | 8.73E-08 | 1.414E-08 | 1.457E-12 | 7.071E-09 |
| Indeno(1,2,3-cd)pyrene | 1.35E-04 | 1.63E-04 | 1.80E-05 | 3.43E-11 | 9.59E-08 | 1.555E-08 | 1.602E-12 | 7.773E-09 |
| Dibenz(a,h)anthracene | 1.78E-05 | 1.80E-05 | 1.80E-05 | ND | ND | ND | ND | ND |
| Benzo(g,h,i)perylene | 5.17E-04 | 5.75E-04 | 1.80E-05 | 1.25E-10 | 3.50E-07 | 5.675E-08 | 5.848E-12 | 2.838E-08 |
| Dioxins and Furans | | | | | | | | |
| 2378-TCDD | 3.26E-09 | 2.54E-09 | 1.03E-09 | ND | ND | ND | ND | ND |
| 12378-PECDD | 6.18E-09 | 5.53E-09 | 1.34E-09 | ND | ND | ND | ND | ND |
| 123478-HXCDD | 3.15E-09 | 2.21E-09 | 6.54E-10 | ND | ND | ND | ND | ND |
| 123678-HXCDD | 3.43E-09 | 2.43E-09 | 7.15E-10 | ND | ND | ND | ND | ND |
| 123789-HXCDD | 5.56E-09 | 3.92E-09 | 1.16E-09 | ND | ND | ND | ND | ND |
| 1234678-HPCDD | 5.55E-09 | 3.19E-09 | 1.88E-09 | 6.61E-16 | 1.85E-12 | 3.000E-13 | 3.092E-17 | 1.500E-13 |
| OCDD | 2.86E-08 | 1.59E-08 | 1.41E-08 | 2.39E-15 | 6.70E-12 | 1.086E-12 | 1.119E-16 | 5.431E-13 |
| 2378-TCDF | 3.91E-09 | 3.17E-09 | 9.42E-10 | ND | ND | ND | ND | ND |
| 12378-PECDF | 4.59E-09 | 3.49E-09 | 1.05E-09 | ND | ND | ND | ND | ND |
| 23478-PECDF | 1.31E-08 | 1.26E-08 | 1.89E-09 | ND | ND | ND | ND | ND |
| 123478-HXCDF | 4.51E-09 | 3.58E-09 | 9.21E-10 | ND | ND | ND | ND | ND |
| 123678-HXCDF | 4.72E-09 | 3.51E-09 | 9.74E-10 | ND | ND | ND | ND | ND |
| 123789-HXCDF | 1.51E-09 | 9.79E-10 | 6.51E-10 | ND | ND | ND | ND | ND |
| 234678-HXCDF | 1.63E-09 | 1.02E-09 | 3.57E-10 | ND | ND | ND | ND | ND |
| 1234678-HPCDF | 2.56E-09 | 1.89E-09 | 9.30E-10 | 2.54E-16 | 7.12E-13 | 1.154E-13 | 1.189E-17 | 5.768E-14 |
| 1234789-HPCDF | 3.32E-09 | 1.96E-09 | 7.90E-10 | ND | ND | ND | ND | ND |
| OCDF | 1.11E-08 | 1.53E-09 | 9.70E-10 | 1.35E-15 | 3.79E-12 | 6.138E-13 | 6.325E-17 | 3.069E-13 |
| Energetics | | | | | | | | |
| Nitrobenzene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 2-Nitrotoluene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 3-Nitrotoluene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 4-Nitrotoluene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| Nitroglycerine | 2.60E-02 | 3.31E-02 | NA | 6.79E-09 | 1.90E-05 | 3.079E-06 | 3.173E-10 | 1.539E-06 |

Table B-1: Air Modeling Output Data for the Long Rifle .22 Caliber Ball Cartridge

| Substance | Trial #1B Measured Actual Concentration (mg/m ³) | Trial #2B Measured Actual Concentration (mg/m ³) | Daily Measured Background Concentration (mg/m ³) | Average Adjusted Emission Factor (lb./item) EF | Average Adjusted Emission Factor (lb./lb. NEW) | Total Mass of Substance Emitted (grams/item) M | Average Modeled Concentration for One Item (grams/m ³) CONC | Substance Emission Rate For One Item (g/item)/sec ER ₁ |
|-----------------------------|--|--|--|--|--|--|---|---|
| 1,3-Dinitrobenzene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 2,6-Dinitrotoluene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 2,4-Dinitrotoluene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 1,3,5-Trinitrobenzene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 2,4,6-Trinitrotoluene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| RDX | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 4-Amino-2,6-Dinitrotoluene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| 2-Amino-4,6-Dinitrotoluene | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| Tetryl | 3.47E-03 | 3.49E-03 | NA | ND | ND | ND | ND | ND |
| HMX | 6.94E-03 | 6.98E-03 | NA | ND | ND | ND | ND | ND |
| Pentaerythritoltetranitrate | 6.94E-03 | 6.98E-03 | NA | ND | ND | ND | ND | ND |
| Dibutyl phthalate | 8.68E-02 | 8.72E-02 | NA | ND | ND | ND | ND | ND |
| Diethyl phthalate | 8.68E-02 | 8.72E-02 | NA | ND | ND | ND | ND | ND |
| Diphenylamine | 8.68E-02 | 8.72E-02 | NA | ND | ND | ND | ND | ND |

Footnotes:

¹ATC = Aberdeen Test Center (for additional information on the data, refer to the Firing Point Emission Study)

NA = Not Applicable

ND = Not Detected

APPENDIX C

HEALTH-BASED SCREENING LEVELS AND ACUTE TOXICITY VALUES

Appendix C: Health-Based Screening Levels and Acute Toxicity Values

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|---------------------------|----------|----------|----------------------|----------|----------------------|-------|------------------------------|------|----------|----------|----------|---------------|
| 1,1,1,2-Tetrachloroethane | 630-20-6 | 2.60E-01 | c | 2.41E-01 | c | | 2.41E-01 | c | 5.15E+04 | | 5.15E+04 | T |
| 1,1,1,1-Trichloroethane | 71-55-6 | 1.04E+03 | nc | 2.30E+03 | nc | | 1.04E+03 | nc | 1.91E+06 | 1.25E+06 | 1.25E+06 | A |
| 1,1,2,2-Tetrachloroethane | 79-34-5 | 3.31E-02 | c | 3.13E-02 | c | | 3.13E-02 | c | 2.06E+04 | | 2.06E+04 | T |
| 1,1,2-Trichloroethane | 79-00-5 | 1.20E-01 | c | 1.12E-01 | c | | 1.12E-01 | c | 1.64E+05 | | 1.64E+05 | T |
| 1,1-Dichloroethane | 75-34-3 | 5.21E+02 | nc | 5.11E+02 | nc | | 5.11E+02 | nc | 1.21E+06 | | 1.21E+06 | T |
| 1,1-Dichloroethene | 75-35-4 | | nc | 3.58E-02 | c | | 3.58E-02 | c | 7.92E+04 | | 7.92E+04 | T |
| 1,2,3-Trichloropropane | 96-18-4 | 9.61E-04 | c | 3.13E-03 | c | | 9.61E-04 | c | 6.03E+04 | | 6.03E+04 | T |
| 1,2,4-Trichlorobenzene | 120-82-1 | 2.08E+02 | nc | 2.08E+02 | nc | | 2.08E+02 | nc | 3.71E+04 | | 3.71E+04 | T |
| 1,2,4-Trimethylbenzene | 95-63-6 | 6.21E+00 | nc | 6.21E+00 | nc | | 6.21E+00 | nc | 1.80E+05 | | 1.80E+05 | T |
| 1,2-Dibromoethane | 106-93-4 | 8.73E-03 | c | 8.24E-03 | c | | 8.24E-03 | c | 1.54E+05 | | 1.54E+05 | T |
| 1,2-Dichlorobenzene | 95-50-1 | 2.09E+02 | nc | 3.29E+02 | nc | | 2.09E+02 | nc | 3.01E+05 | | 3.01E+05 | T |
| 1,2-Dichloroethane | 107-06-2 | 7.39E-02 | c | 6.88E-02 | c | | 6.88E-02 | c | 8.08E+03 | | 8.08E+03 | T |
| 1,2-Dichloropropane | 78-87-5 | 9.89E-02 | c | 9.21E-02 | c | | 9.21E-02 | c | 5.08E+05 | | 5.08E+05 | T |
| 1,3,5-Trimethylbenzene | 108-67-8 | 6.21E+00 | nc | 6.21E+00 | nc | | 6.21E+00 | nc | 3.68E+05 | | 3.68E+05 | T |
| 1,3,5-Trinitrobenzene | 99-35-4 | 1.10E+02 | nc | 1.10E+02 | nc | | 1.10E+02 | nc | 3.00E+04 | | 3.00E+04 | T |
| 1,3-Butadiene | 106-99-0 | 3.74E-03 | c | 3.48E-03 | c | | 3.48E-03 | c | 2.20E+04 | | 2.20E+04 | E |
| 1,3-Dichlorobenzene | 541-73-1 | 3.29E+00 | nc | 1.10E+02 | nc | | 3.29E+00 | nc | 3.61E+04 | | 3.61E+04 | T |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|-----------------------|------------|----------|----------------------|----------|----------------------|-------|----------|----------------------|------|----------|------|----------|---------------|
| 1,3-Dinitrobenzene | 99-65-0 | 3.65E-01 | nc | 3.65E-01 | nc | | 3.65E-01 | nc | | 3.00E+03 | | 3.00E+03 | T |
| 1,4-Dichlorobenzene | 106-46-7 | 3.06E-01 | c | 2.85E-01 | c | | 2.85E-01 | c | | 6.61E+05 | | 6.61E+05 | T |
| 1,4-Dioxane | 123-91-1 | 6.11E-01 | c | 5.69E-01 | c | | 5.69E-01 | c | | 9.00E+04 | | 9.00E+04 | T |
| 1234678-HPCDD | 35822-46-9 | | | | | | | | | | | | |
| 1234678-HPCDF | 67562-39-4 | | | | | | | | | | | | |
| 1234789-HPCDF | 55673-89-7 | | | | | | | | | | | | |
| 123478-HXCDD | 39227-28-6 | | | | | | | | | | | | |
| 123478-HXCDF | 70648-26-9 | | | | | | | | | 8.00E+00 | | 8.00E+00 | T |
| 123678-HXCDD | 57653-85-7 | | | | | | | | | 1.50E+01 | | 1.50E+01 | T |
| 123678-HXCDF | 57117-44-9 | | | | | | | | | 2.00E+00 | | 2.00E+00 | T |
| 123789-HXCDD | 19408-74-3 | 1.48E-06 | c | 1.38E-06 | c | | 1.38E-06 | c | | | | | |
| 123789-HXCDF | 72918-21-9 | | | | | | | | | | | | |
| 12378-PECDD | 40321-76-4 | | | | | | | | | 2.00E+00 | | 2.00E+00 | T |
| 12378-PECDF | 57117-41-6 | | | | | | | | | | | | |
| 1-Butene/Isobutylene | 106-98-9 | | | | | | | | | 6.87E+06 | | 6.87E+06 | T |
| 2,4,5-trichlorophenol | 95-95-4 | 3.65E+02 | nc | 3.65E+02 | nc | | 3.65E+02 | nc | | 3.00E+04 | | 3.00E+04 | T |
| 2,4,6-trichlorophenol | 88-06-2 | 6.20E-01 | c | 6.26E-01 | c | | 6.20E-01 | c | | 3.00E+04 | | 3.00E+04 | T |
| 2,4,6-Trinitrotoluene | 118-96-7 | 2.24E-01 | c | 2.09E-01 | c | | 2.09E-01 | c | | 2.50E+04 | | 2.50E+04 | T |
| 2,4-dichlorophenol | 120-83-2 | 1.10E+01 | nc | 1.10E+01 | nc | | 1.10E+01 | nc | | 3.00E+04 | | 3.00E+04 | T |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|----------------------------|------------|----------|----------------------|----------|----------------------|-------|----------|----------------------|------|----------|------|----------|---------------|
| 2,4-dimethylphenol | 105-67-9 | 7.30E+01 | nc | 7.30E+01 | nc | | 7.30E+01 | nc | | | | | |
| 2,4-dinitrophenol | 51-28-5 | 7.30E+00 | nc | 7.30E+00 | nc | | 7.30E+00 | nc | | 7.50E+03 | | 7.50E+03 | T |
| 2,4-Dinitrotoluene | 121-14-2 | 7.30E+00 | nc | 7.30E+00 | nc | | 7.30E+00 | nc | | 6.00E+02 | | 6.00E+02 | T |
| 2,5-Dimethylbenzaldehyde | 5779-94-2 | | | | | | | | | | | | |
| 2,6-dinitrotoluene | 606-20-2 | 3.65E+00 | nc | 3.65E+00 | nc | | 3.65E+00 | nc | | 6.00E+02 | | 6.00E+02 | T |
| 234678-HXCDF | 60851-34-5 | | | | | | | | | 2.00E+00 | | 2.00E+00 | T |
| 23478-PEGDF | 57117-31-4 | | | | | | | | | | | | |
| 2378-TCDD | 1746-01-6 | 4.48E-08 | c | 4.17E-08 | c | | 4.17E-08 | c | | 4.00E+00 | | 4.00E+00 | T |
| 2378-TCDF | 51207-31-9 | | | | | | | | | 2.00E+00 | | 2.00E+00 | T |
| 2-Amino-4,6-Dinitrotoluene | 35572-78-2 | | | | | | | | | 1.50E+04 | | 1.50E+04 | T |
| 2-Butanone | 78-93-3 | 1.04E+03 | nc | 1.04E+03 | nc | | 1.04E+03 | nc | | 8.85E+05 | | 8.85E+05 | T |
| 2-Butyne | 503-17-3 | | | | | | | | | | | | |
| 2-chloronaphthalene | 91-58-7 | 2.92E+02 | nc | 2.92E+02 | nc | | 2.92E+02 | nc | | 6.00E+02 | | 6.00E+02 | T |
| 2-chlorophenol | 95-57-8 | 1.83E+01 | nc | 1.83E+01 | nc | | 1.83E+01 | nc | | 5.25E+03 | | 5.25E+03 | T |
| 2-Hexanone | 591-78-6 | | | 5.11E+00 | nc | | 5.11E+00 | nc | | 4.09E+04 | | 4.09E+04 | T |
| 2-methylnaphthalene | 91-57-6 | | | 7.30E+01 | nc | | 7.30E+01 | nc | | 2.00E+04 | | 2.00E+04 | T |
| 2-methylphenol | 95-48-7 | 1.83E+02 | nc | 1.83E+02 | nc | | 1.83E+02 | nc | | | | | |
| 2-nitroaniline | 88-74-4 | 2.09E-01 | nc | 2.08E-01 | nc | | 2.08E-01 | nc | | | | | |
| 2-nitrophenol | 88-75-5 | | | | | | | | | | | | |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|----------------------------|------------|----------|----------------------|----------|----------------------|-------|------------------------------|----------|----------|------|----------|---------------|
| 2-Nitrotoluene | 88-72-2 | 3.65E+01 | nc | 3.65E+01 | nc | | 3.65E+01 | nc | | | | |
| 3-Chloropropene | 107-05-1 | 1.04E+00 | nc | | | | | 9.39E+03 | 9.39E+03 | | 9.39E+03 | E |
| 3-nitroaniline | 99-09-2 | | | | | | | | | | | |
| 3-Nitrotoluene | 99-08-1 | 3.65E+01 | nc | 7.30E+01 | nc | | 3.65E+01 | nc | | | | |
| 4,6-dinitro-2-methylphenol | 534-52-1 | | | 3.65E-01 | nc | | 3.65E-01 | nc | 5.00E+02 | | 5.00E+02 | T |
| 4-Amino-2,6-Dinitrotoluene | 19406-51-0 | | | | | | | | | | | |
| 4-bromophenyl-phenylethe | 101-55-3 | | | | | | | | | | | |
| 4-chloro-3-methylphenol | 59-50-7 | | | | | | | | 2.00E+04 | | 2.00E+04 | T |
| 4-chloroaniline | 106-47-8 | 1.46E+01 | nc | 1.46E+01 | nc | | 1.46E+01 | nc | 3.00E+04 | | 3.00E+04 | T |
| 4-chlorophenyl-phenylether | 7005-72-3 | | | | | | | | | | | |
| 4-Ethyltoluene | 622-96-8 | | | | | | | | 1.25E+05 | | 1.25E+05 | T |
| 4-Methyl-2-Pentanone | 108-10-1 | 8.34E+01 | nc | 7.30E+01 | nc | | 7.30E+01 | nc | 3.07E+05 | | 3.07E+05 | T |
| 4-methylphenol | 106-44-5 | 1.83E+01 | nc | 1.83E+01 | nc | | 1.83E+01 | nc | | | | |
| 4-nitroaniline | 100-01-6 | | | | | | | | 9.00E+03 | | 9.00E+03 | T |
| 4-nitrophenol | 100-02-7 | 2.92E+01 | nc | 2.90E+01 | nc | | 2.90E+01 | nc | 3.00E+04 | | 3.00E+04 | T |
| 4-Nitrotoluene | 99-99-0 | 3.65E+01 | nc | 3.65E+01 | nc | | 3.65E+01 | nc | 3.37E+04 | | 3.37E+04 | T |
| acenaphthene | 83-32-9 | 2.19E+02 | nc | 2.19E+02 | nc | | 2.19E+02 | nc | 1.25E+03 | | 1.25E+03 | T |
| Acenaphthylene | 208-96-8 | | | | | | | | 2.00E+02 | | 2.00E+02 | T |
| Acetaldehyde | 75-07-0 | 8.73E-01 | c | 8.13E-01 | c | | 8.13E-01 | c | 1.80E+04 | | 1.80E+04 | E |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|----------------------|-----------|----------|----------------------|----------|----------------------|-------|----------|----------------------|----------|----------|------|----------|---------------|
| Acetone | 67-64-1 | 3.65E+02 | nc | 3.65E+02 | nc | | 3.65E+02 | nc | | 2.37E+06 | | 2.37E+06 | T |
| Acetonitrile | 75-05-8 | 6.20E+01 | nc | 6.21E+01 | nc | | 6.20E+01 | nc | | 1.01E+05 | | 1.01E+05 | T |
| Acetylene | 74-86-2 | | | | | | | | | | | | |
| Acrolein | 107-02-8 | 2.09E-02 | nc | 2.08E-02 | nc | | 2.08E-02 | nc | 2.30E+02 | 2.29E+02 | | 2.30E+02 | E |
| Acrylonitrile | 107-13-1 | 2.83E-02 | c | 2.61E-02 | c | | 2.61E-02 | c | 2.17E+04 | 2.17E+04 | | 2.17E+04 | E |
| Alpha Methyl Styrene | 98-83-9 | 2.56E+02 | nc | 2.56E+02 | nc | | 2.56E+02 | nc | | | | | |
| Aluminum | 7429-90-5 | 5.11E+00 | nc | 3.65E+00 | nc | | 3.65E+00 | nc | | 3.00E+04 | | 3.00E+04 | T |
| Ammonia (NH3) | 7664-41-7 | 1.04E+02 | nc | 1.04E+02 | nc | | 1.04E+02 | nc | 1.75E+04 | 1.75E+04 | | 1.75E+04 | E |
| anthracene | 120-12-7 | 1.10E+03 | nc | 1.10E+03 | nc | | 1.10E+03 | nc | | 6.00E+03 | | 6.00E+03 | T |
| Antimony | 7440-36-0 | | | 1.46E+00 | nc | | 1.46E+00 | nc | | 1.50E+03 | | 1.50E+03 | T |
| Arsenic | 7440-38-2 | 4.47E-04 | c | 4.15E-04 | c | | 4.15E-04 | c | | 3.00E+01 | | 3.00E+01 | T |
| Barium | 7440-39-3 | 5.21E-01 | nc | 5.11E-01 | nc | | 5.11E-01 | nc | | 1.50E+03 | | 1.50E+03 | T |
| Benzaldehyde | 100-52-7 | 3.65E+02 | nc | 3.65E+02 | nc | | 3.65E+02 | nc | | 1.50E+04 | | 1.50E+04 | T |
| Benzene | 71-43-2 | 2.49E-01 | c | 2.16E-01 | c | | 2.16E-01 | c | 1.56E+05 | 1.60E+05 | | 1.56E+05 | E |
| benzo(a)anthracene | 56-55-3 | 2.17E-02 | c | 8.58E-03 | c | | 8.58E-03 | c | | 6.00E+02 | | 6.00E+02 | T |
| benzo(a)pyrene | 50-32-8 | 2.17E-03 | c | 2.02E-03 | c | | 2.02E-03 | c | | 7.50E+03 | | 7.50E+03 | T |
| benzo(b)fluoranthene | 205-99-2 | 2.17E-02 | c | 8.58E-03 | c | | 8.58E-03 | c | | | | | |
| Benzo(e)pyrene | 192-97-2 | | | | | | | | | | | | |
| benzo(g,h,i)perylene | 191-24-2 | | | | | | | | | 3.00E+04 | | 3.00E+04 | T |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS Endpoint | HBSL Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|-----------------------------|-----------|----------|----------------------|----------|----------------------|-------------------|------------------------------|----------|----------|------|----------|---------------|
| benzo(k)fluoranthene | 207-08-9 | 2.17E-01 | c | 8.58E-02 | c | | 8.58E-02 | c | | | | |
| benzyl alcohol | 100-51-6 | 1.10E+03 | nc | 1.10E+03 | nc | | 1.10E+03 | nc | 5.53E+04 | | 5.53E+04 | T |
| Benzyl Chloride | 100-44-7 | 3.96E-02 | c | 3.68E-02 | c | | 3.68E-02 | c | 5.20E+03 | | 5.20E+03 | E |
| Beryllium | 7440-41-7 | 8.00E-04 | c | 7.45E-04 | c | | 7.45E-04 | c | 5.00E+00 | | 5.00E+00 | T |
| bis(2-chloroethoxy)methan | 111-91-1 | | | | | | | | | | | |
| bis(2-chloroethyl)ether | 111-44-4 | 5.82E-03 | c | 5.69E-03 | c | | 5.69E-03 | c | 5.85E+04 | | 5.85E+04 | T |
| bis(2-chloroisopropyl)ether | 108-60-1 | 1.92E-01 | c | 1.79E-01 | c | | 1.79E-01 | c | 6.99E+04 | | 6.99E+04 | T |
| Bis(2-ethylhexyl)phthalate | 117-81-7 | 4.80E-01 | c | 4.47E-01 | c | | 4.47E-01 | c | 1.00E+04 | | 1.00E+04 | T |
| Bromobenzene | 108-86-1 | 1.04E+01 | nc | | | | | | 4.82E+04 | | 4.82E+04 | T |
| Bromodichloromethane | 75-27-4 | 1.08E-01 | c | 1.01E-01 | c | | 1.01E-01 | c | 4.00E+03 | | 4.00E+03 | T |
| Bromoform | 75-25-2 | 1.75E+00 | c | 1.61E+00 | c | | 1.61E+00 | c | 6.20E+03 | | 6.20E+03 | T |
| Bromomethane | 74-83-9 | 5.21E+00 | nc | 5.11E+00 | nc | | 5.11E+00 | nc | 5.82E+04 | | 5.82E+04 | T |
| butylbenzylphthalate | 85-68-7 | 7.30E+02 | nc | 7.30E+02 | nc | | 7.30E+02 | nc | 5.00E+05 | | 5.00E+05 | T |
| Butyraldehyde | 123-72-8 | | | | | | | | 7.38E+04 | | 7.38E+04 | T |
| Cadmium | 7440-43-9 | 1.07E-03 | c | 9.94E-04 | c | | 9.94E-04 | c | 3.00E+01 | | 3.00E+01 | T |
| Calcium | 7440-70-2 | | | | | | | | 3.00E+04 | | 3.00E+04 | T |
| Carbon Dioxide (CO2) | 124-38-9 | | | | | | | | 5.40E+07 | | 5.40E+07 | T |
| Carbon Disulfide | 75-15-0 | 7.30E+02 | nc | 7.30E+02 | nc | | 7.30E+02 | nc | 3.11E+04 | | 3.11E+04 | T |
| Carbon Monoxide (CO) | 630-08-0 | | | | | | 1.00E+04 | 1.00E+04 | 2.30E+05 | | 2.30E+05 | E |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|-------------------------|------------|----------|----------------------|----------|----------------------|-------|----------|----------------------|----------|----------|----------|----------|---------------|
| Carbon Tetrachloride | 56-23-5 | 1.28E-01 | c | 1.18E-01 | c | | 1.18E-01 | c | 1.28E+05 | 1.26E+05 | | 1.28E+05 | E |
| Chlorobenzene | 108-90-7 | 6.21E+01 | nc | 6.21E+01 | nc | | 6.21E+01 | nc | | 1.38E+05 | | 1.38E+05 | T |
| Chlorodifluoromethane | 75-45-6 | 5.11E+04 | nc | 5.11E+04 | nc | | 5.11E+04 | nc | | 4.41E+06 | | 4.41E+06 | T |
| Chloroethane | 75-00-3 | 2.32E+00 | c | 2.16E+00 | c | | 2.16E+00 | c | | 2.64E+06 | | 2.64E+06 | T |
| Chloroform | 67-66-3 | 8.35E-02 | c | 7.73E-02 | c | | 7.73E-02 | c | | 9.76E+03 | | 9.76E+03 | T |
| Chloromethane | 74-87-3 | 1.07E+00 | c | 1.79E+00 | c | | 1.07E+00 | c | | 2.06E+05 | | 2.06E+05 | T |
| Chromium | 7440-47-3 | | c | 1.53E-04 | c | | 1.53E-04 | c | | 1.50E+03 | | 1.50E+03 | T |
| chrysene | 218-01-9 | 2.17E+00 | c | 8.58E-01 | c | | 8.58E-01 | c | | 2.00E+02 | | 2.00E+02 | T |
| cis-1,2-Dichloroethene | 156-59-2 | 3.65E+01 | nc | 3.65E+01 | nc | | 3.65E+01 | nc | | 7.92E+05 | 5.54E+05 | 5.54E+05 | A |
| cis-1,3-Dichloropropene | 10061-01-5 | | | | | | | | | | | | |
| cis-butene | 25167-67-3 | | | | | | | | | 1.72E+04 | | 1.72E+04 | T |
| Cobalt | 7440-48-4 | | | 1.83E-02 | nc | | 1.83E-02 | nc | | 6.00E+01 | | 6.00E+01 | T |
| Copper | 7440-50-8 | | | 1.46E+02 | nc | | 1.46E+02 | nc | | 3.00E+03 | | 3.00E+03 | T |
| Crotonaldehyde | 4170-30-3 | 3.54E-03 | c | | | | | | 5.72E+03 | 5.72E+03 | | 5.72E+03 | E |
| Cumene | 98-82-8 | 4.02E+02 | nc | 4.02E+02 | nc | | 4.02E+02 | nc | | 2.46E+05 | | 2.46E+05 | T |
| dibenz(a,h)anthracene | 53-70-3 | 2.17E-03 | c | 8.58E-04 | c | | 8.58E-04 | c | | 3.00E+04 | | 3.00E+04 | T |
| dibenzofuran | 132-64-9 | 1.46E+01 | nc | 1.46E+01 | nc | | 1.46E+01 | nc | | | | | |
| Dibromochloromethane | 124-48-1 | 8.00E-02 | c | 7.45E-02 | c | | 7.45E-02 | c | | 6.00E+03 | | 6.00E+03 | T |
| Dibromomethane | 74-95-3 | 3.65E+01 | nc | 3.65E+01 | nc | | 3.65E+01 | nc | | 2.50E+05 | | 2.50E+05 | T |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|-------------------------|----------|----------|----------------------|----------|----------------------|-------|----------|----------------------|----------|----------|------|----------|---------------|
| Dibutyl Phthalate | 84-74-2 | 3.65E+02 | nc | 3.65E+02 | nc | | 3.65E+02 | nc | | 1.50E+04 | | 1.50E+04 | T |
| Dichlorodifluoromethane | 75-71-8 | 2.09E+02 | nc | 1.83E+02 | nc | | 1.83E+02 | nc | | 1.48E+07 | | 1.48E+07 | T |
| diethylphthalate | 84-66-2 | 2.92E+03 | nc | 2.92E+03 | nc | | 2.92E+03 | nc | | 1.50E+04 | | 1.50E+04 | T |
| dimethylphthalate | 131-11-3 | 3.65E+04 | nc | 3.65E+04 | nc | | 3.65E+04 | nc | | 1.50E+04 | | 1.50E+04 | T |
| di-n-octylphthalate | 117-84-0 | 7.30E+01 | nc | 7.30E+01 | nc | | 7.30E+01 | nc | | 1.50E+05 | | 1.50E+05 | T |
| Diphenylamine | 122-39-4 | 9.13E+01 | nc | 9.13E+01 | nc | | 9.13E+01 | nc | | 3.00E+04 | | 3.00E+04 | T |
| Ethane | 74-84-0 | | | | | | | | | | | | |
| Ethyl Acetate | 141-78-6 | 3.29E+03 | nc | 3.29E+03 | nc | | 3.29E+03 | nc | | 1.44E+06 | | 1.44E+06 | T |
| Ethyl Acrylate | 140-88-5 | 1.40E-01 | c | | | | | | | 6.14E+04 | | 6.14E+04 | T |
| Ethyl Methacrylate | 97-63-2 | 3.29E+02 | nc | 3.29E+02 | nc | | 3.29E+02 | nc | | | | | |
| Ethylbenzene | 100-41-4 | 1.06E+03 | nc | 1.06E+03 | nc | | 1.06E+03 | nc | | 5.43E+05 | | 5.43E+05 | T |
| Ethylene | 74-85-1 | | | | | | | | | 4.60E+05 | | 4.60E+05 | T |
| fluoranthene | 206-44-0 | 1.46E+02 | nc | 1.46E+02 | nc | | 1.46E+02 | nc | | 3.00E+01 | | 3.00E+01 | T |
| Fluorene | 86-73-7 | 1.46E+02 | nc | 1.46E+02 | nc | | 1.46E+02 | nc | | 7.50E+04 | | 7.50E+04 | T |
| Formaldehyde | 50-00-0 | 1.48E-01 | c | 1.39E-01 | c | | 1.39E-01 | c | 1.23E+03 | 1.23E+03 | | 1.23E+03 | E |
| Freon 113 | 76-13-1 | 3.13E+04 | nc | 3.14E+04 | nc | | 3.13E+04 | nc | | 9.58E+06 | | 9.58E+06 | T |
| Freon 114 | 76-14-2 | | | | | | | | | 2.10E+07 | | 2.10E+07 | T |
| Heptane | 142-82-5 | | | | | | | | | 1.80E+06 | | 1.80E+06 | T |
| hexachlorobenzene | 118-74-1 | 4.18E-03 | c | 3.91E-03 | c | | 3.91E-03 | c | | 7.50E+01 | | 7.50E+01 | T |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|---------------------------|-------------|----------|----------------------|----------|----------------------|----------|----------|----------------------|----------|----------|----------|----------|---------------|
| Hexachlorobutadiene | 87-68-3 | 8.62E-02 | c | 8.03E-02 | c | | 8.03E-02 | c | 3.21E+04 | 3.20E+04 | | 3.21E+04 | E |
| hexachlorocyclopentadiene | 77-47-4 | 7.30E-02 | nc | 7.30E-02 | nc | | 7.30E-02 | nc | | 2.23E+02 | | 2.23E+02 | T |
| Hexachloroethane | 67-72-1 | 4.80E-01 | c | 4.47E-01 | c | | 4.47E-01 | c | | 2.90E+04 | | 2.90E+04 | T |
| Hexaldehyde | 66-25-1 | | | | | | | | | | | | |
| Hexane | 110-54-3 | 2.09E+02 | nc | 2.08E+02 | nc | | 2.08E+02 | nc | | 5.28E+05 | | 5.28E+05 | T |
| HMX | 2691-41-0 | 1.83E+02 | nc | 1.83E+02 | nc | | 1.83E+02 | nc | | | | | |
| Hydrogen bromide | 10035-10-6 | | | | | | | | | 9.93E+03 | | 9.93E+03 | T |
| Hydrogen chloride | 7647-01-0 | 2.08E+01 | nc | 2.08E+01 | nc | | 2.08E+01 | nc | 4.50E+03 | 4.47E+03 | 2.70E+03 | 2.70E+03 | A |
| Hydrogen Cyanide | 74-90-8 | 3.13E+00 | nc | 3.14E+00 | nc | | 3.13E+00 | nc | | 5.17E+03 | | 5.17E+03 | T |
| Hydrogen fluoride | 7664-39-3 | | | | | | | | 1.60E+03 | 1.64E+03 | 1.60E+03 | 1.60E+03 | A |
| indeno(1,2,3-cd)pyrene | 193-39-5 | 2.17E-02 | c | 8.58E-03 | c | | 8.58E-03 | c | | | | | |
| Isobutane | 75-28-5 | | | | | | | | | 9.52E+05 | | 9.52E+05 | T |
| Isooctane | 540-84-1 | | | | | | | | | 3.50E+05 | | 3.50E+05 | T |
| isophorone | 78-59-1 | 7.08E+00 | c | 6.59E+00 | c | | 6.59E+00 | c | | 2.83E+04 | | 2.83E+04 | T |
| Isovaleraldehyde | 590-86-3 | | | | | | | | | | | | |
| Lead | 7439-92-1 | | | | | 2.00E+00 | 2.00E+00 | nc | | 1.50E+02 | | 1.50E+02 | T |
| m/p-Xylene | 108-38-3 10 | 7.30E+02 | nc | 7.30E+03 | nc | | 7.30E+02 | nc | | 6.51E+05 | | 6.51E+05 | T |
| Magnesium | 7439-95-4 | | | | | | | | | 3.00E+04 | | 3.00E+04 | T |
| Manganese | 7439-96-5 | 5.11E-02 | nc | 5.22E-02 | nc | | 5.11E-02 | nc | | 3.00E+03 | | 3.00E+03 | T |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|----------------------------|------------|----------|----------------------|----------|----------------------|----------|----------|----------------------|----------|----------|----------|----------|---------------|
| Methane | 74-82-8 | | | | | | | | | 3.30E+06 | | 3.30E+06 | T |
| Methyl Acrylate | 96-33-3 | 1.10E+02 | nc | 1.10E+02 | nc | | 1.10E+02 | nc | | | | | |
| Methyl Iodide | 74-88-4 | | | | | | | | 1.45E+05 | 1.45E+05 | | 1.45E+05 | E |
| Methyl Methacrylate | 80-62-6 | 7.30E+02 | nc | 7.30E+02 | nc | | 7.30E+02 | nc | | 4.09E+05 | | 4.09E+05 | T |
| Methyl t-Butyl Ether | 1634-04-4 | 3.13E+03 | nc | 3.13E+03 | nc | | 3.13E+03 | nc | | 4.32E+05 | | 4.32E+05 | T |
| Methylene Chloride | 75-09-2 | 4.09E+00 | c | 3.79E+00 | c | | 3.79E+00 | c | 6.96E+05 | 6.94E+05 | | 6.96E+05 | E |
| naphthalene | 91-20-3 | 3.13E+00 | nc | 3.29E+00 | nc | | 3.13E+00 | nc | | 7.86E+04 | | 7.86E+04 | T |
| Nickel | 7440-02-0 | | | 7.30E+01 | nc | | 7.30E+01 | nc | | 3.00E+03 | | 3.00E+03 | T |
| Nitric Acid | 7697-37-2 | | | | | | | | | 2.58E+03 | 1.30E+03 | 1.30E+03 | A |
| Nitrobenzene | 98-95-3 | 2.09E+00 | nc | 2.19E+00 | nc | | 2.09E+00 | nc | | 1.51E+04 | | 1.51E+04 | T |
| Nitroglycerine | 55-63-0 | 4.80E-01 | c | 4.47E-01 | c | | 4.47E-01 | c | | | | | |
| n-nitrosodimethylamine | 62-75-9 | 1.37E-04 | c | 1.23E-04 | c | | 1.23E-04 | c | | 2.50E+03 | | 2.50E+03 | T |
| n-nitroso-di-n-propylamine | 621-64-7 | 9.61E-04 | c | 8.94E-04 | c | | 8.94E-04 | c | | 2.00E+02 | | 2.00E+02 | T |
| n-nitrosodiphenylamine(1) | 86-30-6 | 1.37E+00 | c | 1.28E+00 | c | | 1.28E+00 | c | | | | | |
| o,m,p-Tolualdehyde | 1334-78-7 | | | | | | | | | | | | |
| OCDD | 3268-87-9 | | | | | | | | | 1.50E+02 | | 1.50E+02 | T |
| OCDF | 39001-02-0 | | | | | | | | | 3.00E+02 | | 3.00E+02 | T |
| Octane | 111-65-9 | | | | | | | | | | | | |
| Oxides of Nitrogen (NOx) | 10102-43-9 | | | 3.65E+02 | nc | 1.00E+02 | 1.00E+02 | nc | | 3.08E+04 | | 3.08E+04 | T |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|-----------------------------------|-----------|----------|----------------------|----------|----------------------|----------|----------|----------------------|----------|----------|------|----------|---------------|
| o-Xylene | 95-47-6 | 7.30E+02 | nc | 7.30E+03 | nc | | 7.30E+02 | nc | | 6.51E+05 | | 6.51E+05 | T |
| Particulate Cyanide | 57-12-5 | | | 7.30E+01 | nc | | 7.30E+01 | nc | | 5.00E+03 | | 5.00E+03 | T |
| Particulate Matter <10 micr PM10 | | | nc | | | 5.00E+01 | 5.00E+01 | nc | | | | | |
| Particulate Matter <2.5 mic PM2.5 | | | nc | | | 1.50E+01 | 1.50E+01 | nc | | | | | |
| pentachlorophenol | 87-86-5 | 5.60E-02 | c | 5.22E-02 | c | | 5.22E-02 | c | | 1.50E+03 | | 1.50E+03 | T |
| Pentaerythritoltetranitrate | 78-11-5 | | | | | | | | | 5.00E+01 | | 5.00E+01 | T |
| Pentane | 109-66-0 | | | | | | | | | 1.80E+06 | | 1.80E+06 | T |
| phenanthrene | 85-01-8 | | | | | | | | | 2.00E+03 | | 2.00E+03 | T |
| phenol | 108-95-2 | 2.19E+03 | nc | 2.19E+03 | nc | | 2.19E+03 | nc | | 3.85E+04 | | 3.85E+04 | T |
| Phosphoric acid | 7664-38-2 | 1.04E+01 | nc | 1.06E+01 | nc | | 1.04E+01 | nc | | 3.00E+03 | | 3.00E+03 | T |
| Propane | 74-98-6 | | | | | | | | | 3.78E+06 | | 3.78E+06 | T |
| Propionaldehyde | 123-38-6 | | | | | | | | | 7.50E+04 | | 7.50E+04 | T |
| Propylene | 115-07-1 | | | | | | | | | | | | |
| Propyne | 74-99-7 | | | | | | | | | 2.79E+06 | | 2.79E+06 | T |
| pyrene | 129-00-0 | 1.10E+02 | nc | 1.10E+02 | nc | | 1.10E+02 | nc | | 1.50E+04 | | 1.50E+04 | T |
| RDX | 121-82-4 | 6.11E-02 | c | 5.69E-02 | c | | 5.69E-02 | c | | | | | |
| Selenium | 7782-49-2 | | | 1.83E+01 | nc | | 1.83E+01 | nc | | 6.00E+02 | | 6.00E+02 | T |
| Silver | 7440-22-4 | | | 1.83E+01 | nc | | 1.83E+01 | nc | | 3.00E+02 | | 3.00E+02 | T |
| Styrene | 100-42-5 | 1.06E+03 | nc | 1.04E+03 | nc | | 1.04E+03 | nc | 2.13E+05 | 2.13E+05 | | 2.13E+05 | E |

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| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|----------------------------|------------|----------|----------------------|----------|----------------------|----------|----------|----------------------|----------|----------|----------|----------|---------------|
| Sulfur Dioxide (SO2) | 7446-09-5 | | | | | 8.00E+01 | 8.00E+01 | nc | 7.89E+02 | 7.86E+02 | | 7.89E+02 | E |
| Sulfuric Acid | 7664-93-9 | | | | | | | | 2.00E+03 | 2.00E+03 | | 2.00E+03 | E |
| tert-Butyl Alcohol | 75-65-0 | | | | | | | | | 4.55E+05 | | 4.55E+05 | T |
| Tetrachloroethene | 127-18-4 | 3.31E+00 | c | 3.13E+00 | c | | 3.13E+00 | c | | 6.78E+05 | | 6.78E+05 | T |
| Tetryl | 479-45-8 | 3.65E+01 | nc | 3.65E+01 | nc | | 3.65E+01 | nc | | | | | |
| Thallium | 7440-28-0 | | | 2.56E-01 | nc | | 2.56E-01 | nc | | 3.00E+02 | | 3.00E+02 | T |
| Toluene | 108-88-3 | 4.02E+02 | nc | 4.16E+02 | nc | | 4.02E+02 | nc | 1.88E+05 | 1.89E+05 | | 1.88E+05 | E |
| Total Suspended Particulat | 12789-66-1 | | | | | 5.00E+01 | 5.00E+01 | nc | | | | | |
| trans-1,2-Dichloroethene | 156-60-5 | 7.30E+01 | nc | 7.30E+01 | nc | | 7.30E+01 | nc | 4.95E+04 | 4.95E+04 | 1.11E+06 | 1.11E+06 | A |
| trans-1,3-Dichloropropene | 10061-02-6 | | | | | | | | | | | | |
| Trichlorofluoromethane | 75-69-4 | 7.30E+02 | nc | 7.30E+02 | nc | | 7.30E+02 | nc | | 2.81E+06 | | 2.81E+06 | T |
| Valeraldehyde | 110-62-3 | | | | | | | | | | | | |
| Vanadium | 7440-62-2 | | | 2.56E+01 | nc | | 2.56E+01 | nc | | 1.50E+02 | | 1.50E+02 | T |
| Vinyl Acetate | 108-05-4 | 2.09E+02 | nc | 2.08E+02 | nc | | 2.08E+02 | nc | 1.92E+04 | 1.76E+04 | | 1.92E+04 | E |
| Vinyl Chloride | 75-01-4 | 2.17E-01 | c | 7.20E-02 | c | | 7.20E-02 | c | | 1.28E+04 | | 1.28E+04 | T |
| Zinc | 7440-66-6 | | | 1.10E+03 | nc | | 1.10E+03 | nc | | 3.00E+04 | | 3.00E+04 | T |

| Substance | CAS# | PRG | Toxicity Endpoint | RBC | Toxicity Endpoint | NAAQS | HBSL | Toxicity Endpoint | ERPG | TEEL | AEGL | ATV | ATV Source |
|-----------|------|-----|----------------------|-----|----------------------|-------|------|----------------------|------|------|------|-----|---------------|
|-----------|------|-----|----------------------|-----|----------------------|-------|------|----------------------|------|------|------|-----|---------------|

CAS# = Chemical Abstract Service Number
PRG = Preliminary Remediation Goal (µg/m³)
RBC = Risk-Based Concentration (µg/m³)
NAAQS = National Ambient Air Quality Standard (µg/m³)
HBSL = health-based screening level (µg/m³)
ERPG (E) = Emergency Response Planning Guideline (µg/m³)
TEEL (T) = Temporary Emergency Exposure Limit (µg/m³)
AEGL (A) = Acute Exposure Guideline Level (µg/m³)
c = carcinogen
nc = noncarcinogen

APPENDIX D

RISK ASSESSMENT DATA

Table D-1: Comparison of Modeled Air Concentrations with Health-Based Values

| Long Rifle .22 Caliber Ball Cartridge | | Modeled Distance (meters) 100 | | | DODIC: A106 | | |
|---------------------------------------|------------|-------------------------------|----------|-----------|--------------------------|----------|-----------------------------------|
| Substance* | CAS# | CONC | Cchronic | Cchr/HBSL | Cchr/HBSL>1? | Cacute | Cact/ATV Cact/ATV>1? |
| 1,1,1-Trichloroethane | 71-55-6 | 4.05E-11 | 1.68E-06 | 1.61E-09 | <input type="checkbox"/> | 4.50E-04 | 3.60E-10 <input type="checkbox"/> |
| 1,2-Dichloroethane | 107-06-2 | 4.33E-11 | 7.72E-07 | 1.12E-05 | <input type="checkbox"/> | 1.93E-03 | 2.38E-07 <input type="checkbox"/> |
| 1,3-Butadiene | 106-99-0 | 5.58E-10 | 9.93E-06 | 2.85E-03 | <input type="checkbox"/> | 6.20E-03 | 2.82E-07 <input type="checkbox"/> |
| 1234678-HPCCDD | 35822-46-9 | 3.09E-17 | 1.28E-12 | | <input type="checkbox"/> | 3.44E-10 | <input type="checkbox"/> |
| 1234678-HPCDF | 67562-39-4 | 1.19E-17 | 4.94E-13 | | <input type="checkbox"/> | 1.32E-10 | <input type="checkbox"/> |
| 2-Butanone | 78-93-3 | 3.00E-11 | 1.25E-06 | 1.19E-09 | <input type="checkbox"/> | 1.33E-03 | 1.51E-09 <input type="checkbox"/> |
| Acenaphthene | 83-32-9 | 8.55E-13 | 3.55E-08 | 1.62E-10 | <input type="checkbox"/> | 3.80E-05 | 3.04E-08 <input type="checkbox"/> |
| Acenaphthylene | 208-96-8 | 1.79E-11 | 7.42E-07 | | <input type="checkbox"/> | 7.93E-04 | 3.97E-06 <input type="checkbox"/> |
| Acetaldehyde | 75-07-0 | 1.21E-09 | 2.16E-05 | 2.65E-05 | <input type="checkbox"/> | 1.35E-02 | 7.48E-07 <input type="checkbox"/> |
| Acetone | 67-64-1 | 3.11E-11 | 1.29E-06 | 3.54E-09 | <input type="checkbox"/> | 1.38E-03 | 5.83E-10 <input type="checkbox"/> |
| Acetonitrile | 75-05-8 | 4.67E-10 | 1.94E-05 | 3.13E-07 | <input type="checkbox"/> | 2.08E-02 | 2.06E-07 <input type="checkbox"/> |
| Acetylene | 74-86-2 | 5.51E-09 | 2.29E-04 | | <input type="checkbox"/> | 6.12E-02 | <input type="checkbox"/> |
| Acrolein | 107-02-8 | 1.12E-09 | 4.64E-05 | 2.23E-03 | <input type="checkbox"/> | 1.24E-02 | 5.40E-05 <input type="checkbox"/> |
| Acrylonitrile | 107-13-1 | 3.25E-10 | 5.79E-06 | 2.22E-04 | <input type="checkbox"/> | 3.62E-03 | 1.67E-07 <input type="checkbox"/> |
| Anthracene | 120-12-7 | 4.41E-13 | 1.83E-08 | 1.67E-11 | <input type="checkbox"/> | 1.96E-05 | 3.27E-09 <input type="checkbox"/> |
| Antimony | 7440-36-0 | 4.06E-10 | 1.69E-05 | 1.16E-05 | <input type="checkbox"/> | 1.81E-02 | 1.20E-05 <input type="checkbox"/> |
| Benzene | 71-43-2 | 2.76E-09 | 4.91E-05 | 2.27E-04 | <input type="checkbox"/> | 3.07E-02 | 1.97E-07 <input type="checkbox"/> |
| Benzo(a)anthracene | 56-55-3 | 3.65E-13 | 6.49E-09 | 7.56E-07 | <input type="checkbox"/> | 1.62E-05 | 2.70E-08 <input type="checkbox"/> |

| Long Rifle .22 Caliber Ball Cartridge | | | Modeled Distance (meters) 100 | | | DODIC: A106 | | |
|---------------------------------------|-----------|----------|-------------------------------|-----------|--------------------------|-------------|----------|--------------------------|
| Substance* | CAS# | CONC | Cchronic | Cchr/HBSL | Cchr/HBSL>1? | Cacute | Cact/ATV | Cact/ATV>1? |
| Benzo(a)pyrene | 50-32-8 | 1.46E-12 | 2.59E-08 | 1.28E-05 | <input type="checkbox"/> | 6.48E-05 | 8.64E-09 | <input type="checkbox"/> |
| Benzo(b)fluoranthene | 205-99-2 | 8.15E-13 | 1.45E-08 | 1.69E-06 | <input type="checkbox"/> | 9.06E-06 | | <input type="checkbox"/> |
| Benzo(e)pyrene | 192-97-2 | 1.49E-12 | 6.21E-08 | | <input type="checkbox"/> | 1.66E-05 | | <input type="checkbox"/> |
| Benzo(g,h,i)perylene | 191-24-2 | 5.85E-12 | 2.43E-07 | | <input type="checkbox"/> | 2.60E-04 | 8.66E-09 | <input type="checkbox"/> |
| Benzo(k)fluoranthene | 207-08-9 | 6.71E-13 | 1.20E-08 | 1.39E-07 | <input type="checkbox"/> | 7.46E-06 | | <input type="checkbox"/> |
| Bis(2-ethylhexyl)phthalate | 117-81-7 | 2.45E-10 | 4.36E-06 | 9.76E-06 | <input type="checkbox"/> | 1.09E-02 | 1.09E-06 | <input type="checkbox"/> |
| Butyraldehyde | 123-72-8 | 1.58E-09 | 6.56E-05 | | <input type="checkbox"/> | 7.02E-02 | 9.51E-07 | <input type="checkbox"/> |
| Carbon Dioxide (CO2) | 124-38-9 | 3.45E-06 | 1.43E-01 | | <input type="checkbox"/> | 1.53E+02 | 2.84E-06 | <input type="checkbox"/> |
| Carbon Monoxide (CO) | 530-08-0 | 3.69E-06 | 1.53E-01 | 1.53E-05 | <input type="checkbox"/> | 4.10E+01 | 1.78E-04 | <input type="checkbox"/> |
| Chloromethane | 74-87-3 | 3.60E-13 | 6.41E-09 | 5.99E-09 | <input type="checkbox"/> | 1.60E-05 | 7.77E-11 | <input type="checkbox"/> |
| Chrysene | 218-01-9 | 4.23E-13 | 7.52E-09 | 8.77E-09 | <input type="checkbox"/> | 1.88E-05 | 9.39E-08 | <input type="checkbox"/> |
| Copper | 7440-50-8 | 3.10E-10 | 1.29E-05 | 8.81E-08 | <input type="checkbox"/> | 1.38E-02 | 4.59E-06 | <input type="checkbox"/> |
| Dibutyl phthalate | 84-74-2 | 1.40E-10 | 5.81E-06 | 1.59E-08 | <input type="checkbox"/> | 6.21E-03 | 4.14E-07 | <input type="checkbox"/> |
| Ethane | 74-84-0 | 1.51E-09 | 6.28E-05 | | <input type="checkbox"/> | 1.68E-02 | | <input type="checkbox"/> |
| Ethylene | 74-85-1 | 1.79E-08 | 7.44E-04 | | <input type="checkbox"/> | 7.96E-01 | 1.73E-06 | <input type="checkbox"/> |
| Fluoranthene | 206-44-0 | 5.55E-13 | 2.31E-08 | 1.58E-10 | <input type="checkbox"/> | 2.47E-05 | 8.23E-07 | <input type="checkbox"/> |
| Fluorene | 86-73-7 | 2.10E-12 | 8.72E-08 | 5.98E-10 | <input type="checkbox"/> | 9.33E-05 | 1.24E-09 | <input type="checkbox"/> |
| Formaldehyde | 50-00-0 | 4.01E-09 | 7.13E-05 | 5.13E-04 | <input type="checkbox"/> | 4.45E-02 | 3.62E-05 | <input type="checkbox"/> |
| Hydrogen Cyanide | 74-90-8 | 3.81E-09 | 1.58E-04 | 5.05E-05 | <input type="checkbox"/> | 1.69E-01 | 3.27E-05 | <input type="checkbox"/> |

| Long Rifle .22 Caliber Ball Cartridge | | | Modeled Distance (meters) 100 | | | | DODIC: A106 | | | |
|---------------------------------------|------------|----------|-------------------------------|-----------|--------------------------|----------|-------------|--------------------------|--|--|
| Substance * | CAS# | CONC | Cchronic | Cchr/HBSL | Cchr/HBSL>1? | Cacute | Cact/ATV | Cact/ATV>1? | | |
| Indeno(1,2,3-cd)pyrene | 193-39-5 | 1.60E-12 | 2.85E-08 | 3.32E-06 | <input type="checkbox"/> | 1.78E-05 | | <input type="checkbox"/> | | |
| Lead | 7439-92-1 | 8.52E-08 | 3.54E-03 | 1.77E-03 | <input type="checkbox"/> | 3.79E+00 | 2.52E-02 | <input type="checkbox"/> | | |
| Methane | 74-82-8 | 2.59E-08 | 1.07E-03 | | <input type="checkbox"/> | 1.15E+00 | 3.48E-07 | <input type="checkbox"/> | | |
| Methyl t-Butyl Ether | 1634-04-4 | 1.32E-12 | 5.47E-08 | 1.75E-11 | <input type="checkbox"/> | 5.85E-05 | 1.35E-10 | <input type="checkbox"/> | | |
| Methylene Chloride | 75-09-2 | 8.40E-09 | 1.49E-04 | 3.94E-05 | <input type="checkbox"/> | 9.33E-02 | 1.34E-07 | <input type="checkbox"/> | | |
| Naphthalene | 91-20-3 | 1.75E-10 | 7.29E-06 | 2.33E-06 | <input type="checkbox"/> | 7.80E-03 | 9.92E-08 | <input type="checkbox"/> | | |
| Nitric Acid | 7697-37-2 | 4.07E-09 | 1.69E-04 | | <input type="checkbox"/> | 4.52E-02 | 3.48E-05 | <input type="checkbox"/> | | |
| Nitroglycerine | 55-63-0 | 3.17E-10 | 5.65E-06 | 1.26E-05 | <input type="checkbox"/> | 3.53E-03 | | <input type="checkbox"/> | | |
| OCDD | 3268-87-9 | 1.12E-16 | 4.65E-12 | | <input type="checkbox"/> | 4.97E-09 | 3.32E-11 | <input type="checkbox"/> | | |
| OCDF | 39001-02-0 | 6.33E-17 | 2.63E-12 | | <input type="checkbox"/> | 2.81E-09 | 9.37E-12 | <input type="checkbox"/> | | |
| Oxides of Nitrogen (NOx) | 10102-43-9 | 1.50E-07 | 6.25E-03 | 6.25E-05 | <input type="checkbox"/> | 6.68E+00 | 2.17E-04 | <input type="checkbox"/> | | |
| Particulate Matter <10 micron | PM10 | 1.54E-07 | 6.40E-03 | 1.28E-04 | <input type="checkbox"/> | 1.71E+00 | | <input type="checkbox"/> | | |
| Particulate Matter <2.5 micron | PM2.5 | 1.18E-07 | 4.89E-03 | 3.26E-04 | <input type="checkbox"/> | 1.31E+00 | | <input type="checkbox"/> | | |
| Pentane | 109-66-0 | 2.76E-12 | 1.14E-07 | | <input type="checkbox"/> | 1.22E-04 | 6.81E-11 | <input type="checkbox"/> | | |
| Phenanthrene | 85-01-8 | 2.51E-12 | 1.04E-07 | | <input type="checkbox"/> | 1.11E-04 | 5.57E-08 | <input type="checkbox"/> | | |
| Propylene | 115-07-1 | 3.22E-09 | 1.34E-04 | | <input type="checkbox"/> | 3.58E-02 | | <input type="checkbox"/> | | |
| Pyrene | 129-00-0 | 1.06E-12 | 4.41E-08 | 4.01E-10 | <input type="checkbox"/> | 4.72E-05 | 3.15E-09 | <input type="checkbox"/> | | |
| Styrene | 100-42-5 | 1.37E-10 | 5.69E-06 | 5.45E-09 | <input type="checkbox"/> | 1.52E-03 | 7.14E-09 | <input type="checkbox"/> | | |
| Toluene | 108-88-3 | 2.81E-10 | 1.17E-05 | 2.90E-08 | <input type="checkbox"/> | 3.12E-03 | 1.66E-08 | <input type="checkbox"/> | | |

Monday, August 06, 2001

Long Rifle .22 Caliber Ball Cartridge Modeled Distance (meters) 100 DODIC: A106

| Substance* | CAS# | CONC | Cchronic | Cchr/HBSL | Cchr/HBSL > 1? | Cacute | Cact/ATV | Cact/ATV > 1? |
|-----------------------------|------------|----------|----------|-----------|--------------------------|----------|----------|--------------------------|
| Total Suspended Particulate | 12789-66-1 | 1.52E-07 | 6.30E-03 | 1.26E-04 | <input type="checkbox"/> | 1.69E+00 | | <input type="checkbox"/> |
| Trichlorofluoromethane | 75-69-4 | 1.70E-12 | 7.05E-08 | 9.66E-11 | <input type="checkbox"/> | 7.55E-05 | 2.69E-11 | <input type="checkbox"/> |
| Zinc | 7440-66-6 | 7.15E-10 | 2.97E-05 | 2.71E-08 | <input type="checkbox"/> | 3.18E-02 | 1.06E-06 | <input type="checkbox"/> |

* = Only substances detected in the Firing Point Emission Study are presented in this Appendix. In situations where the substance was detected using more than one sampling method, the higher concentration was used in the risk assessment to maintain a conservative approach.

DODIC = Department of Defense Identification Code

CAS# = Chemical Abstract Service Number

CONC = average modeled concentration for one cartridge (g/m³)

Cchronic = chronic time-averaged concentration (µg/m³)

HBSL = chronic health-based screening level (µg/m³)

Cacute = acute time-averaged concentration (µg/m³)

ATV = acute toxicity value (µg/m³)

Table D-2: Comparison of Modeled Air Concentrations with Health-Based Values: Total Petroleum Hydrocarbons

| Long Rifle .22 Caliber Ball Cartridge | Modeled Distance (meters) | | | DODIC: A106 | |
|---------------------------------------|---------------------------|---------------------------|--------------------------|--------------------------|-------------------------|
| Substance* | CAS# | Cchronic Aliphatic:<=8 | Cchronic Aliphatic:>8 | Cchronic Aromatic:<=8 | Cchronic Aromatic:>8 |
| 1,3-Butadiene | 106-99-0 | 9.93E-06 | | | |
| Acenaphthene | 83-32-9 | | | | 3.55E-08 |
| Acenaphthylene | 208-96-8 | | | | 7.42E-07 |
| Anthracene | 120-12-7 | | | | 1.83E-08 |
| Benzene | 71-43-2 | | | 4.91E-05 | |
| Benzo(a)anthracene | 56-55-3 | | | | 6.49E-09 |
| Benzo(a)pyrene | 50-32-8 | | | | 2.59E-08 |
| Benzo(b)fluoranthene | 205-99-2 | | | | 1.45E-08 |
| Benzo(e)pyrene | 192-97-2 | | | | 6.21E-08 |
| Benzo(g,h,i)perylene | 191-24-2 | | | | 2.43E-07 |
| Benzo(k)fluoranthene | 207-08-9 | | | | 1.20E-08 |
| Chrysene | 218-01-9 | | | | 7.52E-09 |
| Fluoranthene | 206-44-0 | | | | 2.31E-08 |
| Fluorene | 86-73-7 | | | | 8.72E-08 |
| Indeno(1,2,3-cd)pyrene | 193-39-5 | | | | 2.85E-08 |
| Naphthalene | 91-20-3 | | | | 7.29E-06 |
| Pentane | 109-66-0 | | 1.14E-07 | | |

| Long Rifle .22 Caliber Ball Cartridge | | Modeled Distance (meters) | | 100 | DODIC: A106 | |
|--|----------|---------------------------|--------------------------|--------------------------|-------------------------|--|
| Substance* | CAS# | Cchronic Aliphatic:<=8 | Cchronic Aliphatic:>8 | Cchronic Aromatic:<=8 | Cchronic Aromatic:>8 | |
| Phenanthrene | 85-01-8 | | | | 1.04E-07 | |
| Propylene | 115-07-1 | 1.34E-04 | | | | |
| Pyrene | 129-00-0 | | | | 4.41E-08 | |
| Styrene | 100-42-5 | | | | 5.69E-06 | |
| Toluene | 108-88-3 | | | 1.17E-05 | | |
| Total (µg/m³) | | 1.44E-04 | | 6.08E-05 | 1.44E-05 | |
| Derived Health-Based Screening Level (µg/m³) | | 1.92E+04 | 1.04E+03 | 4.17E+02 | 2.09E+02 | |
| Cchronic/HBSL | | 7.49E-09 | | 1.46E-07 | 6.90E-08 | |
| Is this ratio >1? | | No | No | No | No | |

* = Only substances detected in the Firing Point Emission Study are presented in this Appendix. In situations where the substance was detected using more than one sampling method, the higher concentration was used in the risk assessment to maintain a conservative approach.

DODIC = Department of Defence Identification Code
CAS# = Chemical Abstract Service Number
Cchronic = chronic time-averaged concentration (µg/m³)
HBSL = chronic health-based screening level (µg/m³)

APPENDIX E

FACT SHEET SUBMITTED TO THE
U.S. ARMY ENVIRONMENTAL CENTER

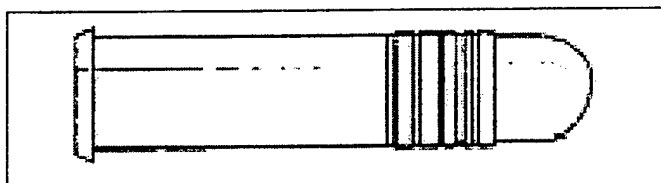
U.S. Army Environmental Center

Training Munitions Fact Sheet

Long Rifle .22 Caliber Ball Cartridge

Department of Defense Identification Code: A106

Breathing air emissions from the Long Rifle .22 Caliber Ball Cartridge will not impact the health of residents who live near Army training facilities.



To be fully prepared to protect our country, U.S. soldiers must train with many different weapons and munitions, including the Long Rifle .22 Caliber Ball Cartridge (.22 Caliber Ball). This training is important because it helps prepare our soldiers for a variety of combat situations. While the Army recognizes the value of such comprehensive training on our installations, we also work hard to ensure the safety and health of surrounding communities.

WILL BREATHING AIR EMISSIONS FROM THE LONG RIFLE .22 CALIBER BALL CARTRIDGE AFFECT MY HEALTH?

To answer this question, the U.S. Army tested the air emissions that are released when the .22 Caliber Ball is fired. The information gathered during these tests was then analyzed to determine if there would be a potential for health effects from inhalation to residents who live near training areas. Study results, generated using conservative methods, showed that offsite residents breathing air as close as 100 meters (328 feet or about the length of a football field) from the firing location are safe from these emissions. At most locations, training areas are at least 1,000 meters (over half a mile) away from populated areas and the distance to firing locations may be even farther.

HOW WAS THE STUDY CONDUCTED?

To gather data for this study, the .22 Caliber Ball was fired from a rifle in a test chamber. The air in the chamber was then tested to identify the types and amounts of substances released. About 200 different substances were looked for during this part of the study.

This information was then used in an U.S. Environmental Protection Agency (USEPA) approved air model (a computer program that allows estimation of air concentrations) to determine the amount of each substance to which someone

living near a training site might be exposed. Downwind concentrations were estimated based on a typical use scenario for the .22 Caliber Ball during training exercises. Since this study did not look at any one specific training area, the assumptions used in the model would, in most cases, predict higher downwind air concentrations than those expected at an actual training site.

These estimated air concentrations were then compared to screening levels established by the USEPA and other federal agencies. If the air concentrations are below these screening levels, they are considered safe for the general population, including sensitive people such as the sick, elderly, and children.

WHAT ARE THE STUDY LIMITATIONS?

Many steps were taken to ensure that the results of this study are protective of residents who live near training facilities. However, as with any study, this study has limitations. For example, the study does not consider exposure to other types of munitions that could also be used during the same training event. Due to these limitations, conservative model conditions were used to ensure the protection of public health from breathing .22 Caliber Ball air emissions.

WHAT EXACTLY IS THE LONG RIFLE .22 CALIBER BALL CARTRIDGE?

The .22 Caliber Ball is a type of ball ammunition used for marksmanship practice and match use. The .22 Caliber Ball consists of a copper alloy cartridge case and a lead-antimony bullet. It also contains a propelling charge that consists mostly of nitrocellulose. Nitrocellulose is the primary ingredient in smokeless propellant (for both military and commercial use) and is also used in the production of lacquers and artificial leathers. Each .22 Caliber Ball is about as long as the width of a quarter.

WHERE CAN I GET MORE INFORMATION?

For more information on the .22 Caliber Ball or other military munitions, please call the Army Environmental Hotline at 1-800-USA-3845, visit our Web site at www.aec.army.mil, or e-mail t2hotline@aec.apgea.army.mil.